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(54) Title: MODEL FOR TESTING IMMUNOGENICITY OF PEPTIDES

(57) Abstract

Assay methods for determining whether a peptide is likely to be immunogenic are based on a computer modeling of binding to a Class II MHC DR1 receptor. This is confirmed by competitive inhibition binding assays. The peptides are useful for eliciting an immune response for vaccination or the production of antibodies or T-cells.

Applicants: Alexander Gad et al.

Serial No.: 09/816,989 Filed: March 23, 2001

Exhibit 11

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MODEL FOR TESTING IMMUNOGENICITY OF PEPTIDES

Government Interest

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The invention described herein may be manufactured, licensed and used by or for governmental purposes without the payment of any royalties to us thereon.

Cross Reference

This application is a continuation-in-part of U.S. Patent application Serial No. 08/064,559, filed May 21, 1993, and the present application incorporates U.S. Patent Application Serial No. 08/064,559 in its entirety by reference.

15 Field of the Invention:

This invention relates to a means of predicting potential of a peptide for eliciting immune response.

Background of the Invention:

Among the numerous steps required for an immunological response to occur is the presentation of the antigen by macrophages to the B-cell or T-cell. This presentation is mediated by the Class I and Class II major histocompatibility complex (MHC) molecules on the surface of the cell. The MHC molecules hold antigens in the form of the peptide fragments and together with the receptor molecule on the T-cells, form a macromolecular complex that induces a response in the T-cell. Therefore, a necessary step in an immune response is the binding of the antigen to the MHC.

Recent single crystal X-ray structures of human and murine Class I MHC's have been reported. Analysis of these crystal structures have shown that antigenic peptides lie in the so-called binding cleft for presentation to the T-cell. This cleft is formed by α_1 and α_2 domains and by β -strands from each domain forming the floor. Furthermore, the sequence polymorphism among Class I molecules can result in alterations of the surface of the cleft forming different pockets. Peptide side chains may insert into these pockets. Thus, different pockets may interact with different side chains. This implies the mechanism for the peptide specificity of Class I MHC's. Peptides bound to the Class I MHC's in the crystal structures were found to have both the amino and carboxy termini tightly held by the MHC. There were few interactions near the middle of the cleft. Hence the bound peptide is allowed to bend slightly in the center. observed binding mode helped to explain the apparent partial specificity of peptide sequence and the allowed variation in peptide length found among peptides isolated from Class I

The precise mode of binding of peptides to Class II MHC molecules is less clear. While a single crystal X-ray diffraction structure for the HLA-DR1 MHC has been shown, the coordinates have remained unavailable. However, currently available theoretical and experimental results help form a hypothesis that the binding of a peptide to Class II MHC is similar to that observed with Class I. First, it is noted that the Class II binding cleft is structurally similar to

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that of Class I. This was concluded based upon a sequence analysis of 26 Class I and 54 Class II amino acid sequences.

Unlike with Class I molecules, self-peptides isolated from murine I-Ab and I-Eb, from murine I-Ad and from human HLA-DR1 molecules were found to be varied in size (13 to 25 residues long). The peptides isolated from the murine I-Ab and I-Eb molecules had heterogenous carboxy termini while those from I-Ad and HLA-DR1 had ragged termini at both ends. The varying lengths indicate that the amino and carboxy termini of the peptides were not critical for the binding. One or both termini may protrude from the binding site and be available for further processing. The residues critical for binding were proposed to be at the ends of the peptide as opposed to the center.

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Summary of the Invention:

It is the purpose of this invention to provide a method for preliminary screening of peptides for ability to elicit an immune response. Structural homology techniques were used to model a receptor (the Class II MHC is exemplified). This model makes it possible to preliminarily screen peptides for antigenic properties. By modifying the peptide to "fit" into the receptor it is possible to identify methods of rendering non-immunogenic peptides immunogenic.

The preliminary screening of peptides for immunogenicity comprises the steps of (1) creating a molecular model of a receptor followed by minimizing the model created, 2) modeling a peptide to be tested and minimizing the model of the peptide, then testing the

fit of the model of the peptide into the model of the receptor to produce a composite minimized receptor/minimized peptide model. Upon finding an acceptable fit, the peptide may then be screened by a binding assay for actual binding to Class II MHC as a further tes for immunogenicity.

It has been found that when the model of the peptide can not be fitted into the model of the receptor, the peptide will lack immunogenicity. While not all peptide models which can be made to "fit" into to model of the receptor will be effective as immunogens the screening methods of the invention may make it possible to avoi undue biological testing of inappropriate peptides. By using the model, it is also possible to alter peptides to accommodate the receptor. Hence, the invention has both predictive and drug design applications.

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Brief Description of the Figures:

Fig. 1 shows the HLA-aw68 $lpha_1$ and $lpha_2$ domains with DR1 $lpha_1$ and $oldsymbol{eta}_1$

Figs. 2-30 are a printout of the minimized coordinates of the receptor.

Figs. 31 and 32 shows the effects of various peptides inhibiting the binding of labeled hemagglutinin in a competitive binding assay.

25 Detailed Description of the Invention:

In order to understand and better predict peptide interaction with Class II MHC's and as an aid for synthetic peptide vaccine design, a structural homology model of HLA-DR1 molecule was made

using the Class I HLA-aw68 as a reference molecule. For purposes of this analysis, numerous conserved residues were aligned leading to a proposed three-dimensional model for the Class II structure very similar to that of Class I. This model retained the overall conformation of a Class I MHC and agreed with a considerable amount of the published data. Furthermore, peptides shown to bind to DR1 were docked in the binding cleft of the model and analyzed. The results agree with the experimental binding data presented here. Hence, it is shown that the structural homology model reported here is useful for screening Class II MHC functionality.

It had been hypothesized that few peptide residues may be required for binding to DR1. By substituting residues into the influenza hemagglutinin 307-319 T-cell epitope (HA) it had been determined that a single tyrosine at 308 was required for binding. A synthetic peptide with the tyrosine at position 308 and a lysine at 315 was found to bind DR1 as well as the native peptide. Hence, it was concluded that few peptide residues determine the high affinity binding to DR1.

The peptides produced according to the present invention may be used alone or chemically bound to another peptide and/or carrier in order to elicit an immune response. An immune response is elicited by administering a peptide to an animal in an effective dose and by an effective route of administration. Typically the peptide will be administered with an immunologically acceptable carrier. The routes of administration, dosages, times between multiple administrations will be based on the particular peptide and are standard operations of those skilled in the art.

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Of particular interest are peptides from pathogenic microorganisms and neoplasms. In such an example, a vaccine may be formed with the peptide and any known immunological carrier and may be administered prophylactically or therapeutically. The immune response may be elicited for a number of reasons other than for prophylaxis or therapy such as increasing antibody production for the harvesting of antibodies, or increasing specific B-cell or T-cell concentration for the production of hybridomas or cellular therapy.

The choice of host animals is limited only to those capable of an immune response. Preferred hosts are mammals, more preferred are humans.

The vaccine may contain plural peptides with each peptide corresponding to the same or different antigens. The peptides may be used unbound or they may be chemically bound to another peptide or an unrelated protein or other molecule. A preferred vaccine preparation contains a plurality of peptides chemically bound to a larger more immunogenic peptide.

The peptide may be adsorbed, bound or encapsulated in a biodegradeable microsphere, microcapsule, larger carrier or a combination of these. The carrier may have a slow or controlled release property thereby releasing the peptide under appropriate conditions and times for enhanced immunization. This is particularly important when administering the peptide orally where stomach acid can degrade the peptide.

Another embodiment of the present invention is to modify the amino acid sequence of a peptide to enhance its immunogenicity.

This is done by modifying the natural peptide sequence to bind to

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the Class II MHC receptor DR1 with superior binding affinity for a Class II MHC receptor DR1 than the natural peptide sequence. This modified peptide is considered a synthetic peptide. Alternatively, the sequence may be modified to have a greater inhibition of HA (306-318) binding to a Class II MHC receptor DR1.

Many amino acid changes are acceptable in the formation of a synthetic peptide. The changes may be for similar types of amino acids such as leucine for isoleucine or they may be for diverse types such as tyrosine for lysine.

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Materials and Methods:

The structural homology model for the DR1 Class II MHC was constructed using the QUANTA molecular modeling package (vision 3.2, Molecular Simulations, Inc., Burlington, MA) with the CHARMM and Protein Design modules. After alignment of the sequences as described below, gaps and loops were energy minimized using 100 steps of steepest descents minimization followed by 100 steps of adopted basis set Newton-Rapheson (ABNR) minimization. were closed using a fragment database from a selected set of high-resolution crystal structures. The resulting structure was minimized in vacuo using 1000 steps of steepest descents followed by an additional 1000 steps of ABNR minimization. A distance related electrostatic function was used in all calculations with a dielectric constant of 1.0. Non-bound parameter lists were updated every 20 steps with a cutoff distance of 15.0Å. Non-bonded calculations were performed using a shifted potential function between 11.0A and 14.0A. An extended atom set was used with only

polar hydrogen atoms specifically placed. There were no explicit hydrogen bond energy calculations performed.

All peptides were initially modeled using QUANTA in an extended chain conformation and subjected to 500 steps of ABNR minimization. The resulting structures remained essentially in extended chain conformations. Individual peptides were manually docked in several different orientations into the binding cleft region of the minimized DR1 structure. The resulting bimolecular complex was subjected to 5000 steps of steepest descents minimization with non-bonded interactions updated every five steps. After minimization, bound peptides remained essentially in extended chain conformations. The lowest energy complexes for each peptide were selected for further analysis.

The selected peptide and DR1 complexes and the minimized DR1 model were subjected to the following molecular dynamics regimen: 300 steps of heating to 300°K, 600 steps of equilibration at 300°K, and 1100 steps of production dynamics. During this simulation, the DR1 C α atoms were constrained in their starting positions. All non-bonded interaction parameters were as stated for the minimization procedure. The lowest energy structure during the course of the production dynamics was selected and subjected to the 5000 step minimization procedure described previously with the C α restraints removed. The resulting structures were used for the binding energy calculations and for hydrogen bonding analysis.

Hydrogen bonds were determined using the QUANTA default parameters. Maximum allowed distances were 2.5Å between a hydrogen and the acceptor atom and 3.3Å between the donor and acceptor atoms

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The minimum angle allowed between any set of atoms forming a hydrogen bond was 90°.

Competitive Inhibition Binding Assay:

HA peptide (the influenza hemagglutinin 307-319 T-cell epitope) was labeled with ¹²⁵I. The labeled HA peptides were then allowed to interact with purified DR1 molecules during incubation to allow formation of peptide/DR1 complexes. After incubation, the peptide/DR1 composition was exposed to a native gel for chromatographic separation or passed through a spun column to separate labeled peptide/DR1 complex and free labelled peptide. When unlabeled peptides were added before incubation of labeled HA peptides and DR1, and if the unlabelled peptides had capacity for binding to DR1 simultaneous with ¹²⁵I-HA, there was a resultant decrease in radioactive signal associated with the DR1. The extent of this decrease directly related to the binding capacity of the unlabeled unknown peptide.

Structural Homology Model for the DR1 Molecule:

The structural homology model was created, the reference molecule being the crystal structure of HLA-aw68. The HLA-aw68 coordinates and subsequent sequence were obtained from the entry 2HLA in the Brookhaven Protein Data Bank released January 15, 1991, which is incorporated herein by reference. The sequence for the DR1 molecule was for the α_1 domain was reported by Klein and for the β_1 domain, the study reported by Todd et al. (Nature 329, 599 (1987)).

The sequence alignment is based on Brown et al. (Nature 332, 845 (1988)). The complete alignment and numbering scheme for both

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are seen in Figure 1. The Class II, eta_1 and Class I $lpha_2$ domains regions were conserved with some variations at the ends where the two MHC's have different loop regions. The fourth B-strand in the α_1 domain of HLA-aw68 (residues 30-38) is disrupted in the DR1 model. Only three residues are in a β -sheet conformation, probably due to the inserted glycine at position 28 before the strand and the large deletion in the loop region immediately after the strand. two alpha-helical regions are clearly maintained. Both helices have been observed to be discontinuous in the Class I molecules and are similar in the DR1 model. The α_i domain helix is long and curves from residues 49α to 76α without significant disruption. essentially a single continuous helix. However, the α_2 helical region is broken into two separate helices as with the Class I molecules. A short helix (52-63) is separated from a longer helix (68-94) by a deformed region without secondary structure. This deformation is more pronounced in the DR1 model as opposed to the Class I molecules due to an insertion.

Influenza Hemagglutinin Peptide with DR1:

The amino acid residues 307-319 of influenza hemagglutinin (Pro-Lys- Tyr-Val-Lys-Gln-Asn-Thr-Leu-Lys-Leu-Ala-Thr) make up a well-documented linear T-cell epitope which has been shown to be HLA-DR1 restricted. With the demonstration that the influenza hemagglutinin epitope (referred to as the HA peptide) binds DR1, it was chosen to be modeled into the binding cleft.

The peptide was initially inserted into the cleft so that Leu 11 HA was in the vicinity of the hydrophobic pocket. This allowed Asn 7 to be near the middle charged and polar groups of the cleft.

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The remaining residue of the motif (Lys 2) was near the vicinity of the remaining charged and polar residues at the end of the cleft.

The only adjustment to the starting conformation was a slight rearrangement of the terminal peptide proline and Tyr 3 to alleviate obvious bad contacts.

After the energy minimization of the bimolecular complex, the total energy was reduced to 483 kcal/mol. This reduction in energy was accomplished by alleviation of several bad contacts and also be formation of several hydrogen bonds. The sticking feature of this mode is lack of hydrogen bonds in the carboxy terminal half of the peptide. Only one hydrogen bond is identified between the backbone carbonyl group of Leu 9 and the side chain of the β_1 Asn 77. In contrast, the amino terminal half has eleven identified interactions. Four of these interaction involve the peptide backbone residues Tyr 3, Val 4, and Gln 6. The remainder involve the side chains of Lys 2, Tyr 3, Lys 5 and Gln 6. Interestingly, Lys 5 is involved in more interactions (three) than Lys 2 (only 2). No interactions were observed as anticipated with Asn 7. Instead, it was the glutamine at position 6 donating a hydrogen bond to the $lpha_1$ Asn 62. No interactions were observed for the amino and carboxy termini.

HA-YK Peptide with DR1:

The binding of the HA-YK peptide (Ala-Ala-Tyr-Ala-Ala-Ala-Ala-Ala-Ala-Ala-Ala-Lys-Ala-Ala) to the DR1 model was tested. In aligning the peptide in the cleft, it was deemed logical to insert the tyrosine residue into the hydrophobic region of the binding cleft. The lysine would then be in position to interact with the

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nydrophilic groups in the other half of the cleft. The resulting peptide orientation is the opposite that used for the HA and the CS3 (defined below) peptides. With the peptide oriented as described, the final docking position for the peptide was unclear. The hydrophobic pocket is quite large, and, at least in this model, could accommodate the peptide tyrosine in a number of positions by sliding the peptide lengthwise through the cleft. However, repositioning the peptide also repositions the lysine. There were primarily two positions for the lysine: one with the lysine inside the cleft and the second with it outside. Of the two positions, the former was the lower in energy by 46 kcal/mol and had the greater number of interactions with the protein (11 vs. 7). Thus, the preferred orientation of the peptide appears to be with the lysine inside the binding cleft region.

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CS3 subunit Pilin Peptide with DR1:

The suspected T-cell epitope for CS3 pilus subunit 63-78

(Ser-Lys-Asn-Gly-Thr-Val-Thr-Trp-Ala-His-Glu-Thr-Asn-Asn-Ser-Ala)

was modeled with the DR1 molecule. The peptide was inserted with

lysine inside the cleft in the hydrophilic region. This placed the

Thr 5 in the center of the binding cleft and the tryptophane

(residue 8) near the hydrophobic region. The resulting minimized

model had ten interactions between the peptide and the protein,

three interactions with the peptide backbone and five with the

peptide side chains. The remaining two were with the amino terminal

of the peptide. All of the interactions were in either the first

three residues, His 10 or Glu 11 in the peptide. No interactions

were observed in the center of the cleft or residues four through nine.

CFA/1 with DR1:

A peptide identified as CFA/1 (colonization factor antigen)

(Val-Gly-Lys-Asn-Ile-Thr-Val-Thr-Ala-Ser-Val-Asp-Pro) was prepared

and an attempt was made to "fit" the molecule into the cleft of the

DR1. The lysine at position 3 prevented insertion of the peptide.

10 Results:

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The peptides chosen to dock in the DR1 model are shown in Table

1. The peptides were docked manually in several orientations into
the DR1 model. The peptides were then tested in biological binding
assays with the following results:

Table I

Peptide	Molecular Model predicted binding	Binding in the bioassay
HA (influenza hemagglutinin)	Yes	Yes
HA-YK (synthetic peptide)	Yes	Yes
CS3 Pilin subunit	Yes	Yes
CFA/1	No	No

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Quantitative measurement of the inhibition of CS3 63-78 and HA 306-318 as compared to controls is shown in Fig. 31.

The binding energy was calculated as the difference between the final DR1 and peptide complex and the sum of the energies for the minimized DR and peptide models individually. The data is shown in Table II.

Table II.

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Peptide	Protein	Residues	Sequence	Binding Energy (kcal/mol)
АН	Influenza hemagglutinin	306-318	PKYVKQNTLKLAT	-283
на-үк	synthetic peptide		ААҮАААААКАА	-216
CS3	CS3 pilin subunit	63-78	SKNGTVTWAHETNNSA	-245

CS60 and CS6B with DR1

Colonization factor antigen IV (CFA/IV is an antigen on the surface of many enterotoxigenic *E. coli* one component of which is CS6. CS6 has two major subunits and a number of minor subunits. Several peptides from CS6 have been sequenced and assayed for potential inhibition of radiolabeled HA (306-318)/DR1 complex as a measure of immunogenicity. The sequences of the subunits are shown in Table III.

Table III.

TADIE TII.		
Peptide	Amino Acid Residues	Sequence
CS6α6	63-75	DEYGLGRLVNTAD
CS6α7	80-92	IIYQIVDEKGKKK
CS6α8	111-123	LNYTSGEKKISPG
CS6B1	3-15	WQYKSLDVNVNIE
CS6B2	42-54	QLYTVEMTIPAGV
CS6ß3	112-124	TSYTFSAIYTGGE
CS6ß4	123-135	GEYPNSGYSSGTY
CS6ß5	133-145	GTYAGHLTVSFYS

These peptides were assayed for inhibition of radioactively labeled HA(306-318)/DR1. The results are demonstrated in Fig. 32.

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BNSDOCID: <WO___9531997A1_I_>

The foregoing description of the specific embodiments reveal the general nature of the invention so that others can, by applying current knowledge, readily modify and/or adapt for various applications such specific embodiments without departing from the generic concept, and, therefore, such adaptations and modifications should and are intended to be comprehended within the meaning and range of equivalents of the disclosed embodiments. It is to be understood that the phraseology or terminology employed herein is for the purpose of description and not of limitation.

All references mentioned in this application are incorporated by reference.

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We Claim:

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1. A method of preliminarily screening peptides for immunogenicity comprising the steps of:

- 1) creating a molecular model of receptor DR1 Class II MHC and minimizing the model of the DR1;
- 2) modeling a peptide to be tested and minimizing the model of the peptide; and
- 3) testing fit of model obtained in step 2 into the model

 10 obtained in step 1 to produce a composite receptor/peptide model.
 - 2. A computerized model comprising a model of the DR1 molecule having fitted in a cleft therein a model of a peptide.
- 3. A method of claim 1 wherein, additionally, the receptor/peptide model is subjected to computer-simulated heating.
 - 4. A method of claim 1 further comprising, assaying the peptide by competitive inhibition binding to a Class II MHC receptor DR1.
 - 5. A minimized peptide capable of binding to a Class II MHC receptor DR1 and inhibiting the binding of HA (306-318).
- 6. A synthetic peptide, wherein the amino acid sequence of the
 minimized peptide according to claim 5 has been modified to have a
 superior binding affinity for a Class II MHC receptor DR1 to form at
 least a portion of the synthetic peptide.

7. A synthetic peptide, wherein the amino acid sequence of the minimized peptide according to claim 5, has been modified to have greater inhibition of HA (306-318) binding to a Class II MHC receptor DR1 to form at least a portion of the synthetic peptide.

- 8. A synthetic peptide according to claim 6, wherein an amino acid has been modified from a charged amino acid to an uncharged amino acid.
- 9. A synthetic peptide according to claim 7, wherein an amino acid has been modified from a charged amino acid to an uncharged amino acid.
- 10. A synthetic peptide according to claim 8, wherein said uncharged amino acid is alanine.
 - 11. A synthetic peptide according to claim 9, wherein said uncharged amino acid is alanine.
- 12. A minimized peptide according to claim 5, wherein the sequence is selected from the group consisting of PKYVKQNTLKLAT, AAYAAAAAKAA and SKNGTVTWAHETNNSA.
- 13. A minimized peptide according to claim 5, wherein the sequence is contained in a CFA.

14. A minimized peptide according to claim 13, wherein the sequence is selected from the group consisting of DEYGLGRLVNTAD, IIYQIVDEKGKKK, LNYTSGEKKISPG, WQYKSLDVNVNIE, QLYTVEMTIPAGV, TSYTFSAIYTGGE, GEYPNSGYSSGTY and GTYAGHLTVSFYS.

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- 15. A vaccine comprising:
 - a minimized peptide according to claim 5; and an immunologically acceptable carrier.
- 10 16. A vaccine comprising:
 - a synthetic peptide according to claim 6; and an immunologically acceptable carrier.
 - 17. A vaccine comprising:
- a synthetic peptide according to claim 7; and an immunologically acceptable carrier.
- 18. A method of eliciting an immune response in an animal comprising administering said animal with the vaccine according to claim 15.
 - 19. A method of eliciting an immune response in an animal comprising administering said animal with the vaccine according to claim 16.

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20. A method of eliciting an immune response in an animal comprising administering said animal with the vaccine according to claim 17.

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                     -50.39491 -50.76541 101.74483 A1
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         2 LYS
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                                           97.74074 Al
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  59
        6 VAL
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                     -48,60121 -30,33669 101,01420 A1
        5 VAL
                C.3
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FIG. 2

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63	. 6 VAL	CG1	47 07672	-36.30119	101.00126	Al	ų U	17-0-0000
62	6 VAL	CG2	19.10544	-37.46647	100.11057	λl	6	0.00000
€3	6 VAL	c	-49.19435	-35.00190	103.02675	Al	6	0.00000
64	6 VAL	ō	-50,2101B	-34.34041	102.87457	X1	6	0.00000
65	7 ILE	N	-48,11527	-34.52120	103.65431	A1	7	0.00000
66	7 ILE	н	-47.35972	-35.13148	103.91409	X1	7	0.00000 0.00000
67	7 ILE	СЪ	-48.09506	-33.08697	103.98819	λl	7	0.00000
68	7 ILE	CB	-48.69197		105.39701	Υï	7	0.00000
69	7 ILE	CG2	-47.96322	-33.71317	106.43001		7	0.00000
. 70	7 ILE	CG1	-48.74316	-31.39586	105.81727		7	0.00000
71	7 ILE	CD	-49.28846	-31.20489	107.23523		7	0.00000
72	7 ILE	С	-46.69381	-32.50114	103.87753		, ל	0.00000
73	7 ILE	0	-45.72315	-33.10109	104.32366	VT	8	0.00000
74	8 ILE	И	-46.61414	-31.32789	103.23109	VI	8	0.00000
75	8 ILE	н		-30.79777	103.00074 102.75879		8	0.00000
76	8 ILE	CA		-30.85113			8	0.00000
77	8 ILE	CB	-45.18096	-31.21426	100.47371		8	0.0000
78	9 ILE	CG2	-46.47692	-30.98133	100.51961	A1	8	0.0000
79	8 ILE	CG1	-44.01581	-30.33370	101.12422	Al	8	0.0000
80	8 ILE .		-42.65515	-29.37643	103.03239	λl	8	0.0000
81	8 ILE	C		-28.43922			8	0.00000
82	8 ILE		-45.69216	-20.43322	103.70229		9	0.0000
83	9 GLN	Ν .	-43.86371	-29.19466 -29.97409	104.05676	Al	9	0.0000
84	9 GTN	H	-43.33/1/ -43.27910	-23.37903	103.71549		9	0.0000
85	9 GIN	CA.	-43.27910	-27.43231	105.14730		9	0.00000
25	9 GLN	CB	44.97213	-27.13894	105.94550	Al	9	0.00000
87	a Cin a Ciń	CG CD	-43 92932	-26.69043	107.36359	Al	9	0.0000
85 89	9 GLN 9 GLN	OE1	-44.59992	-27.05224	108.31811	Al	9	0.00000
90	9 GLN	NE2	-42.89278	-25.86874	107.50418	X1	9	0.00000
91	9 GLN	HE21	-42.31986	-25.58069	106.73881	A1	9	0.00000
92	9 GLN	HE22	-42.66204	-25.53260	108.41526	Al	9	0.00000 0.00000
93	9 GLŃ	C	-42.00840	-27.79728	102.89330	W.T	9 9	0.00000
94	9 GLN	Ο.	-41.07030	-28.56703	103.06902		10	0.00000
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96	10 ALA	H		-26.24022	101.81084		3.0	0.00000
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98	10 ALA	CB		-26.95993	101.25674		10	0.00000
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101	ii Gru	И	-39,11637 -38.51800	-24.95043	100.86462	Al	11	0.00000
102	ij Gru	H		-23.61077		Al	11	0.00000
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104 105	11 GLU 11 GLU	CG	-38.29740	-24.15021		Al	11	0.00000
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111	12 PHE	N	-37.86688	-21.87280	99.78525	λl	12	0.00000
112	12 PHE	Н	-38.38856	-21.25141	100.37827	Al	12.	0.00000
113	12 PHE	CA	-37.20863	-21.33691	98.59813	Y.I	12 12	0.00000
114	12 PHE	CB	-38.26225		97.51950 96.27668	V 1	12	0.00000
115	12 PHE	CG		-21.86124	96.27665		12	0.0000
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122	12 PHE	0	-35 16677	-20.22034	99.15293	A l	13	0.00000
123 124	13 TYR 13 TYR	H H	-34 71888	-21.10856	99.03348	F. 1	1.3	0.00000
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        19 SER
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189	19 SER	OG	34.3706	4 -11.8772	7 103.10841	λ1	19	0.00000
190					9 103.40254		19	0.00000
3 92	19 SER	. C			6 100.02197		19	0.00000
152	19 SER	0	-36.4951	8 -12.8640	99.00144	Al	19	0.00000
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156	20 GLY	С		7 -15.8974			20	0.00000
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198	21 GLU	N	The second secon	•	2 100.25477		21	0.00000
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200 201	21 GLU 21 GLU	CA CB			9 100.27114 0 98.79559		21	0.00000
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211	22 PHE	CB		-20.31207			22	0.00000
212 213	22 PHE	CG		-21.48591			22	0.00000
213	22 PHE 22 PHE	CD1			5 105.44915		22	0.00000
215	22 PHE 22 PHE	CD2 CE1			5 104.14562 3 106.34948		22 · 22	0.00000
216	22 PHE	CE2			105.04232		22	0.00000
217	22 PHE	CZ			106.14281		22	0.00000
218	22 PHE	C.			101.58028		22	0.00000
219	22 PHE	0			101.55111		22	0.00000
220	23 MET	N			101.32298		23	0.00000
221	23 MET	н			101.30500		23	0.00000
222	23 MET	CA			101.02967		23	0.00000
223	23 MET	CB		-22.73266			23	0.00000
224	23 MET	CG		-23.35273			23	0.00000
225	23 MET	SD		-23.10231			23	0.00000
226	23 MET	CE		-24.67640			23	0.00000
227 228	23 MET 23 MET	0			101.21411		23 23	0.00000
229		N			100.62115		24	0.00000
230	• •	Н		-26.26850			24	0.00000
231		CA			102.69060		24	0.00000
232		CB	-47.06193		103.21139		24	0.00000
233	24 PHE	CG	-46.27878	-23.27691	104.43188	X 1	24	0.00000
234		CD1			105.60532	A1	24	0.00000
235		CD2			104.40480		24	0.00000
236		CEl			106.75356		24	0.00000
237		CE2	-46.89083	-21.60180	105.55271	Al .	24	0.00000
238 239		CZ			106.72434		24	0.00000
240		C.,			101.85776		24 24	0.00000
241		N 0			101.74049			0.00000
242					102.20704			0.00000
243		CA .			100.90165			0.00000
244		23	-48.21052		99.52046			0.0000
245		CG	-49.19634		98.47932 2			0.00000
246	25 ASP. C		-49.75730	-27.46672	97.77589 1			0.0000
247	•	DD 2	-49.39656		98.38197			0.00000
248	25 ASP (101.56183 7			0.00000
245	25 ASP C		-40.57317	-29.35714	102.56241	.1		0.00000
250	26 PHE 1	J	-50.15819	-29.70261	101.00377	1		0.00000
251 252	26 PHE H	١.,	-50.60050 ·	-27.38662	100.15349 7	.1		0.00000 0.00000
232	26 PHE C	: A	-30.30030 ·	-30.77/91	101.21030 /	••	• •	0.0000

FIG. 5

PCT/US94/05697

, · , ,	عادت عاديناهي	ٔ	<u> ಜ್ಞಾ</u>	25 14:58:	68 1993.		5	
253	26 PHE		51.4237	8 -30.8309	3 102.8200	. A1	26	0.00000
254	26 PHI		52.6035	6 -29.9203	8 102.57724	Al	26	0.00000
255	26 PHE		-52.5039	8 -28.5567	4 102.91777) Al	26	0.00000
256	26 PHE	_			5 101.99011		26	0.00000
257	26 PHE				1 102.66706		26	0.00000
258	26 PHE				5 101.74257		26	0.00000
259	26 PHE				7 102.07912		26	0.00000
260	26 PHE				100.51070		26	0.00000
261	26 PHE			7 -31.3821			26	0.00000
262	27 ASP				100.89199		27	0.00000
263	27 ASP				101.83544		27	0.00000
264	27 ASP				3 100.01741 3 100.24733		27 27	0.00000
265 266	27 ASP 27 ASP				100.73899		27	0.00000
267	27 ASP	CG OD1		-36.23164		A1	27	0.00000
268	27 ASP	OD2			100.47314		27	0.00000
269	27 ASP	C		-34.05698			27	0.00000
270	27 ASP	õ		-34.10386			27	0.00000
271	28 GLY	N		-33.92246	4.		28	0.00000
272	28 GLY		· -49.78144				28	0.0000
273	28 GLY	CA	-49.88845	-33.86633	96.98221	A1	28	0.00000
274	28 GLY	C	-50.00090	-32.51821	96.28701	Al	28	0.00000
275	28 GLY	0	-A9_04304	.1-32.03342	95.69731	Al	28	0.00000
276	29 ASP	N	-51.21573	-31.96145	96.33541	Al	29	0.00000
277	29 ASP	H	-51.93955	-32.39453	96.87595	λl	29	0.00000
278	29 ASP	Cλ	-51.52130	-30.85124	95.43198	Al	29	0.00000
279	29 ASP	CB.		-31.44665			29	0.00000
280	29 ASP	CC		-30.54960			29	0.00000
281	29 ASP	OD1		-30.32445	92.44162		29	0.00000
282	29 ASP	002		-30.10545			29	0.00000
283	29 ASP	C		-29.64631	96.07233		29	0.00000
284	29 ASP	0		-28.59407			29 30	0.00000 0.00000
285 286	30 GLU	×		-29.81701	97.32405 97.90251		30	0.00000
287	30 GLU 30 GLU	H		-30.57647 -28.70871	97.86160		30	0.00000
288	30 GLU	CA CB		-29.18152	98.54724	וג	30	0.00000
289	30 GLU	CG		-29.71506	97.67444		30	0.00000
290	30 GLU	CD		-29.50058	98.43828		30	0.00000
291	30 GLU	OE1		-30.47326	98.78969		30	0.00000
292	30 GLU	OE2		-28.34297	98.66447		30	0.00000
293	30 GLU	C		-27.80995	98.82987		30	0.00000
294	30 GLU			-28.20198	99.84476	A1	30	0.00000
295	31 ILE	N		-26.53268	98.46810	λl	31	0.00000
296	31 ILE	. н	-53.28532	-26.29454	97.65353	A1	31	0.00000
297	31 ILE	CX	-51.98442	-25.52556	99.18511		31	0.00000
298	31 ILE	CB	-51.81933		98.23317		31	0.00000
299	31 ILE	CG2	-53.16329		•	Al	31	0.00000
300	31 ILE	CG1	-50.75168		98.71478		31	0.00000
301	31 ILE	CD	-50.41981		97.66271		31	0.00000
302 303	31 ILE	С	-52,51316		100.56211		31 31	0.00000 0.00000
304	31 ILE 32 PHE	14 ·	-53.70233 -51.54696		100.82175		32	. 0.00000
305	32 PHE	H .	-50.59061	•	101.18020		32	0.00000
306	32 PHE	CA	-51.85606		102.84494		32	0.00000
307	32 PHE		-50.76201				32	0.00000
308	32 PHE	CG	-51.19588		104.74893		32	0.00000
309	32 PHE	CD1	-50.23527		105.15120		32	0.00000
310	32 PHE	CD2	-52.49149				32	0.0000
311	32 PHS	CEI	-50.57386	-28.14516	106.12524	Al	32	0.00000
312	32 PHE	CE2	-52.82688	-27.19600	106.29575	A.1	32	0.00000
313	32 PHE	ÇZ	-51.86506	-28,15152	106.69566	VΊ	32	0.00000
314	32 PHE	С	-51.84357	-23.07181	103.07985	¥.1	32	0.00000
315	32 PHE	0	-52.76581	-22.::086	103.62205	λl	32	0.00000
316	33 HIS	N	-50.69098	-22.50451	102.70813	ř. J	33	0.00000

FIG. 6

_نةنار.	يلاندادك وحود	>	The Feb	25 14:58:	48 1993 .		6	
317	33 HIS	н	0.0431	2 -22.9883	4 102.10786	. A1	33	0.00000
318	33 HIS		0.2257	6 -21.2172	7 103.23242		33	0.00000
37.0	SIK EE		-49.6590	g -21.5154	0 104.63303		33	0.00000
320	33 HIS	CG	-49.2298	4 -20.3214	2 105.45366		33	0.00000
321	33 HIS		48.0282	6 -19.7279	3 105.3720		33	0.00000
322	33 HIS		-47.2899	9 -19.9275	7 104.75923		33	0.00000
323	33 XIS		-49.9721	4 -19.6882	B 106.45019		33	0.00000
324	33 HIS			8 -18.70683			33 33	0.00000
325	33 HIS			0 -18.7265 1 -20.7847			33	0.00000
326	33 HIS			5 -21.6120			33	0.00000
327	33 HIS 34 VAL		-48.3372	7 -19.5044	102.33695		34	0.00000
328 329	34 VAL		-49 1942	9 -18.80802	102.90183		34	0.00000
330	34 VAL		-47.5177	6 -19.11490	101.64269		34	0.00000
331	34 VAL		-47.8130	3 -18,42567	7 100.28392	Al	34	0.00000
332	34 VAL			6 -18.41158			34	0.00000
333	34 VAL	CG2		1 -19.06363			34	0.00000
334	34 VAL	С	-46.79652	2 -18.14692	2 102.56509	YJ	34	0.00000
335	34 VAL	0	-47.41849	-17.54298	103.42874	λl	34	0.00000
336	35 ASP	N	-45.47963	-18.03426	102.37666	Al	35 35	0.00000 0.00000
337	35 ASP	H			101.76710 102.93355		35	0.00000
338 .339	35 ASP 35 ASP	CA CB			101.94639		35	0.00000
340	35 ASP	CG			101.88569		35	0.00000
341	35 ASP	OD1	-46 85196	-15.31279	100.93528	A1	35	0.00000
3 42	35 ASP	OD2	-46.43780	-14.28565	102.78747	λl	35	0.00000
343	35 ASP	C			104.41152		35	0.00000
344	35 ASP	0	-45.03304	-17.43137	105.27733	¥1	35	0.0000
345	36 MET	N .	-44.91212	-15.26386	104.68945	λl	36	0.0000
346	36 MET	H.	:	-14.58792			36	0.00000
347	36 MET	CA		-14,74244			36	0.00003
348	36 MET	CB			106.01845		36	0.00000 0.00000
349	36 MET	CC.			105.50403		36 36	0.00000
350	36 MET	SD		-11.71848 -10.40582			36	0.00000
351 352	36 MET	CE C	-45.60204	-16 27022	106.53712	A1	36	0.00000
353	36 MET	0		-14.73983			36	0.0000
354	37 ALA	ĸ			105.58618		37	0.0000
355	37 ALA	Я	-47.21869	-14.87225	104.60519	X1	37	0.00000
356	37 ALA	CA			106.05297	Al	37	0.0000
357	37 ALA	СВ		-13.19637			37	0.00000
358	37 ALA	C	•	-15.61256			37	0.00000
359	37 ALA	0		-16.18787			37	0.00000
360		. N	-50.76933	-15.79391	106.59397		38 38	0.00000
361 362	38 TX2	H CA	-52.00981	-16 48765	106.23832	Al	30	0.00000
363	38 LY6	CB	-52.90628	-16.55867	107.48308	λl	38	0.00000
364	38 LYS	CG	-52.41585	-17.40517	108.60236	Al	38	0.00000
365	30 LYS	CD,	-53.40991	-18.62084	108.86983	Al	38	0.0000
366	38 TJ2	CE:	-53.42547	-19.66155	107.75111	Al	38	0.00000
367	38 LYS	ĸz	-54.78503	-20.17861	107.56038	A1	38	0.00000
368	38 LYS		-54.79226	-20.92377	106.83759	A1	38 38	0.00000
369	38 LYS	HZ2	-55.17879	-20.59510	108.44046	VT	38 - 38	0.00000
370 371	38 LYS	HZ3	-32.44/47	-15 73427	107.27224	- E 1	3 B	0.00000
3/1 372	38 LYS	CO	-52 77571	-14 51200	105.12989	X1	38	0.00000
373	39 LYS	N	-54 35457	-16 44351	104.17702	Al	39	0.00000
374	39 LYS	ĥ v	-53 59774	-15.96113	103.33578	A1	3,9	0.00000
375	39 LYS	CA	-53.67982	-17.86931	104.22041	A.	39	0.00000
376	39 LYS	CB	-55.18971	-18.00506	104.55346	A.1	39	0.00000
377	39 LYS	CG	-56.24681	-17.64030	103.42417	Ä1	39	0.00006
378	39 LYS	CD	-56.27039	-15.52580	102.€2130	λl	39	0.00000
375	39 LYS	CE	-56.64532	-15.69151	101.13213	Al .	39	0.0000
350	39 LYS	ΝI	-55.66137	-17.54092	100.43802	A:	33	0.00000

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381	39 LYS	H21	55.9564	5 -18.5458	1 100.36313		0.00000
382	39 LYS.	HZ2	-55.3986	3 -17.2672	5 99.46103		
3 6 3	39 LYS	HZ3	-54.7248	6 -17.51923	1 100.90690	A1 3	
384	39 LYS	C	-53.4442	3 -18.44283	3 102.84433	A1 3	0.0000
385	39 LYS	0	-53,4335	1 -17.67518	8 101.88936	A1 3	9 0.00000
386	40 GLU	N	-53.4133	2 -19.77420	102.71885	A1 4	0.00000
387	40 GLU	н	-53.1615	4 -20.43693	1 103.42697	A1 4	0.0000
388	40 GLU	CA	-54.0889	2 -20.14645	5 101.48663	A1 4	0.0000
389	40 GLU	CB	-53.2213	2 -20.84643	3 100.44056	A1 4	0.00000
390	40 GLU	CG	-52.68203	3 -19.82757	99.41114	A1 4	0.00000
391	40 GLU	9	-53.7822	7 -18.90107			0.00000
392	40 GLU	OE1	-53 606B	1 -17.68372	98.89245		0.00000
393	40 GLU	OE2	-54 88246	6 -19.35560	98.59153		·
394	40 GLU	C	-55 46629	-20.74566	101.61501		
395	40 GLU	o	-56 (203)	-20.18304	101.08479	A1 40	
. 396	41 THR	N	-55 56371	-21 R4976	102.35754	A1 4:	
_	41 THR	11			102.86321		
397	41 THR	CA	-56 94476	-22.20412	102.30963	A1 41	
398 399	41 THR	CB	-56.65476		102.16109	A1 41	
400	41 THR	OGI	-50.33011	24.05050 -24.75407	101.74066	A1 41	
401					101.33840		
401	41 THR 41 THR	HG1			103.43152		
		CG2					
403	41 THR	С			103.42344		
404	41 THR	0			104.56718		
405	42 VAL	N	-59.12316	-22.51525	103.00455		
406	42 VAL	Н	-59.25687	-22.81422	102.09633		
407	42 VAL	CA	-60.29134	-22.03291	103.80812	A1 42	
408	42 VAL	CB	-61.57611	-22.26846	102.98525	A1 42	
409	42 YAL	CGI	-62.83989	-21.83994	103.74041	A1 42	
410	42 VAL	CG2	-61.49852	-21.55078	101.63610	A1 42	
411	42 VAL	Ç	-60.39368	-22.75550	105.24170	A1 42	
412	42 VAL	0			106.20404	A1 42	
413	43 TRP	N·		-24.07997			
414	43 TRP	н.		-24.53437	104.20145		
415	43 TRP	CÀ		-24.82113			
416	43 TRP	CB.		-26.25321			
417	43 TRP	CG	-60.96314	-26.72686	104.73262	እ 1 43	
418	43 TRP	CD2	-59.88712	-27.12198	103.92127	A1 43	
419	43 TRP	CE2	-60.47947	-27.49629	102.60102	ሕ ጋ 43	
420	43 TRP	CE3	-58.50029		104.12574		
421	43 TRP	CD1			103.93459	ሕ1 43	
422	43 TRP	NEl	-61.84174	-27.31306	102.67766		
423	43 TRP	HE1			101.97106		0.0000
424		CZ2	-59.62565	-27.96652	101.58525		0.0000
425	43 TRP (CŻ3			103.07694		0.0000
426	43 TRP (CH2	-58.24204	-28.07937	101.82597	Al 43	0.00.000
427	43 TRP	c ·	-59.07150	-24.84373	107.16564	Al 43	0.00000
428	43 TRP (0	-58.85050	-25.68073	108.02836	Al 43	0.00000
429	44 ARG 1	N	-58.22471	-23.84563	106.89519	Al 44	0.0000
430	44 ARG I	н	-58.31398	-23.27977	106.07364	A1 44	0.0000
431		CA			107.89144		0.00000
432		CB	-55.83244	-23.69727	107.26013	A1 44	0.0000
433	••	CG	-54.63741		108.22026		0.0000
434	•	CD	-53.30915	-23.64273	107.47006		0.00000
435		NE:	-52 18625	-23.26000	108.32672	A1 44	0.00000
436	•	HE	-52.36221	-22.91543	109.25139	R1 44	0.00000
437		cz	-50.93149		107.86475		0.00000
438	•	\H1	-49.9235B		108.56844		0.00000
439		HH11	-48.98076		108.24673		0.00000
440			-50.05162		109.51721		0.00000
441		vH2	-50.67215		106.69352		0.00000
4 4 2			-49.73581		106.35139		0.00000
443			-51.41769		106.14195		0.00000
444	ARG C		-57.42751		108.38029		0.00000
	11 NV3 C	•	24.52/31			• •	

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445	' 44 AR		£ 530£	4 -21.3920	1 108.88943	E 1	44	0.00000
446	45 LE		-58.6527	0 -21.5568	6 108.18548	λl	45	0.00000
447	45 LEI		-59.3684	6 -22.0638	2 107.69729	አነ	45	0.00000
448	45 LE	J CA	-58.9467	9 -20.2189	7 108.70611	λl	45	0.00000
445	45 LE	CB	59.8726	7 -19.4651	5 107.74942	λl	45	0.00000
450	45 LE		-59:1505	6 -18.9699	0 106.49579	Al	45	0.00000
451	45 LET		-60:1453	4 -18.4190	6 105.47407	λl	45	0.00000
452	45 LEC		-58:1043	1 -17.9228	9 106.87944	λl	45	0.00000
453	45 LEU				4 110.09168		45	0.00000
454					6 110.79199		45 46	0.00000
455	46 GLU				4 110.48954 B 109.87631		46	0.00000
456	46 GLU	•			7 111.86481		46	0.00000
457 458	46 GLU				111.95437		46	0.00000
459	46 GLU				111.51772		46	0.00000
460	46 GLU				111.46594		46	0.00000
461	46 GLU	OE1	-62.26867	-25.47222	110.36424	ÀΊ	46	0.00000
462	46 GLU	OE2			112.52280		46	0.00000
463	46 GLU	С			112.92683		46	0.00000
464	46 GLU	.0			114,10493		46	0.00000
. 465	47 GLU	31			112.44220		47	0.00000
466 -467	47 GLU 47 GLU	H			111.48920		47 47	0.00000
468	47 GLU	CA CB			112.39059		47	0.00000
469	47 GLU	CG			111.78977		47	0.00000
470	47 GLU	CD			112.78701		47	0.00000
471	47 GLU	0E1			112.37878		47	0.00000
472	47 GLU	OE2			113.95212		47	0.00000
473	47 GLU	С		•	113.91510		17	0.00000
474	47 GLU	0	-56.67000	-19.35742	113.26194	A1	47	0.00000
475	48 PHE	N	-56.83807	-20.40716	115.24713	YJ	48	0.0000
476	48 PHE	Р.			115.72305		48	0.00000
477	48 PHE	CY			115.96362		48	0.00000
478	48 PHE	CB			117.41420 117.70829		48 48	0.00000 0.00000
479 480	48 PHE	CG	-50 71127	-19 00320	117.44215	21	48	0.00000
481	48 PHE	CD2	-58.25459	-17.19966	118.24455	A.1	48	0.00000
482	48 PHE	CE1			117.71246		48	0.00000
483	48 PHE	CE2	-59.38940	-16.40286	118.51469	አ1	4 B	0.00000
484	48 PHE	CZ	-60.68202	-16.90878	118.24779	Al	4 B	0.0000
485	48 PHE	C			115.93777		48	0.0000
486	48 PHE	0			115.72587		48	0.00000
487	49 GLY	И			116.19781		49	0.00000 0.00000
488	49 GLY	H C	-56.32069	-16.72101	116.34575	Al al	49 49	0.00000
.489 490	49 GLY	C CV	-59.23420	-16.30320	116.68056	14	49	0.00000
491	49 GLY	0	-51.85815	-16.69886	116.10714	Al	49	0.00000
492	50 ARG	N ·			117.81779		50	0.00000
493	50 ARG	н	-53.93666	-17.67284	118.24069	Al	50	0.00000
494	50 ARG	CX			118.42224		50	0.00000
495	50 ARG	CB.	-52.27437	-18.92760	119.68340	A1	50	0.00000
496	50 ARG	CG	-51.10260	-19.32736	120.58254	Al	50	. 0.00000
497	50 ARG	CD			121.80047		50 50	0.00000 0.00000
498 499	50 ARG	ΝE	-50.3/432	-10.44743	122.64580	7.1 2.1	50	0.00000
500	50 ARG	HE C2	-50 24449	-31 E3184	123.26571	A1	50	0.00000
501	50 ARG	NH1			124.04503		50	0.00000
502	50 ARG		-49.05579	-22.71179	124.52612	P.]	50	0.00000
503	SO ARG	HH12	-48.49360	-21.13233	124.16453	~ 1	50	0.00000
504	50 ARG	NH2	-51.15926	-22.58629	123.10823	ř. 1	50	0.00000
505	50 ARG	HH24	-51.08073	-23.47414	123.56513	ሉ)	50	0.00000
506	50 ARG		-51.95092	-22.42795	122.51852	ĂÌ • ì	50	0.00000
507	50 ARG	c	-51.06703	-19.09561	117.49074	r.i .·	50	0.00000
508	50 ARG	0	-49.84240	-19.05!33	117.52826	r	50	. 0.00000

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509	51 PHE	и	· 1.84915 -19.87778 116.73839	A1 51	0.00000
510	51 PHE	н	2.84564 -19.76657 116.76797		0.00000
513	SI PHE	Cλ	-51.29477 -20.78796 115.73402		0.00000
512	51 PHE	CB	-52.51672 -21.52503 115.15449		0.00000
513	51 PHE	CG	-52.25099 -22.49194 114.02359	A1 51	0.00000
514	51 PHE	CD 1	-51.87888 -23.82851 114.29503	A1 51	0.00000
515	51 PHE	CD 2	-52.45788 -22.06484 112.69289	אז 51	0.00000
516	51 PHE	CEI	-51.72483 -24.74421 113.22969	A1 51	0.00000
517	51 PHE	CE2	-52.30778 -22.98082 111.63059	A1 51	0.00000
518	51 PHE	CZ	-51.94314 -24.31860 111.90000	A1 51	0.00000
519	51 PHE	C	-50.50268 -20.00696 114.69371	A1 51	0.0000
520	51 PHE	0	-49.31281 -20.20801 114.47961		0.00000
521	52 ALA	.κ	-51.20469 -19.03212 114.10202	A1 52	0.00000
522	52 ALA	н	-52.16823 -18.89764 114.33870		0.00000
523	52 ALA	CA	-50.54896 -18.16486 113.11957		0.00000
523 524	52 ALA	CB	-51.52058 -17.09216 112.62511	A1 52	0.00000
525	52 ALA	C	-49.28257 -17.48933 113.62687		0.00000
526	52 ALA	0	-48.27008 -17.39834 112.94510		0.00000
527	53 SER	N	-49.35763 -17.04955 114.88764		0.00000
528	53 SER		·· -50.22411 -17.11498 115.38567		0.00000
529		CA	-48.18100 -16.47055 115.53815		0.0000
530	53 SER 53 SER	CB	-48.58146 -16.02158 116.95383		0.00000
531			-4759593 -15.15541 117.52845		0.00000
532	53 SER 53 SER	0G	-47.85158 -14.90357 118.42105		0.00000
		НG	-46.99433 -17.42737 115.57426		0.00000
533		C	-45.89463 -17.11790 115.12568	A1 53	0.00000
534 535		0	-47.26082 -18.64220 116.08200		0.00000
536	54 PHE 54 PHE); L	-40.18040 -10.87361 116.41568		0.00000
537	•	H	-46.18727 -19.64350 116.09999		0.00000
538	54 PHE 54 PHE	CA	-46.69548 -20.99079 116.63413		0.00000
		CB.	-46.90625 -20.99411 118.13255		0.00000
539		CG	-48.11656 -21.50316 118.65539		0.00000
540	54 PHE	CD1	-45.89246 -20.53119 119.00621		0.00000
541	54 PHE	CD2	-48.31310 -21.55376 120.05266		0.00000
542	54 PHE	CEl	• •		0.00000
543	54 PHE	CE2	-47.30008 -21.09192 120.92307 A		0.00000
544	54 PHE	CZ	-45.57270 -19.90650 114.73758 A		0.00000
545	54 PHE	C	-44.36030 -19.93109 114.55246 A	A1 54	0.00000
546	54 PHE	0	-46.46681 -20.08976 113.76558 A	A1 55	0.00000
547 548	55 GLU 55 GLU		-47.45338 -20.04226 113.95516 A	A1 55	0.00000
•	55 GLU 55 GLU	H.	-45.97322 -20.41545 112.42886 A	A1 55	0.00000
549	55 GLU	CA	-47.14512 -20.81915 111.54876 A		0.00000
550	55 GLU	CB CG	-47.92567 -21.96126 112.19274 P	A1 55	0.00000
551 5 5 2	55 GLU	.CD	-49.04456 -22.36202 111.27677	A1 55	0.00000
553	55 GLU	OE1		A1 55	0.00000
554	55 GLU	OE2	-49.80524 -21.50984 110.81474		0.00000
555	55 GLU	C.	-45.19354 -19.30188 111.76318		0.00000
55.6	55 GLU	ō:	-44.12678 -19.50387 111.19343		0.00000
557	56 ALA	И	-45.73650 -18.08673 111.89532 J		0.00000
558	56 ALA	H	-46.62336 -17.96765 112.34890		0.00000
559	56 ALA	CA	-45.00414 -16.92733 111.38640 3	41 56	0.00000
560		CB	-45.80074 -15.63871 111.59969 J		0.00000
561	56 ALA	c	-43.63772 -16.77849 112.02791 4	11 56	0.00000
562	56 ALA	ō	-42.62065 -16.60634 111.36878 J	1 56	0.00000
563		N	-43.63088 -16.91454 113.35929 A	11 57	0.00000
564		н	-44.48826 -17.03983 113.86832 4		0.00000
565		Cλ	-42.35063 -16.88200 114.06847 A		0.00000
566		CB	-42.61967 -17.03049 115.56930 F		0.00000
567	- • -	CG	-41.40659 -15.75355 116.45895 F		0.00000
566		CD	-41.80070 -16.94589 117.90800 A		0.00000
569		OZ1	-42.29879 -16.06206 118.58910 F		0.00000
570		NE2	-41.55425 -10.16011 116.30366 A		0.00000
571	57 GLN	2221	-41.14709 -18.87176 117.81428 A		0.00000
572	57 GLN	222	-41.78115 -15.36728 119.33313 A	12 57	0.00000
	J. J			-	

./pR1_	ಟ್ಯಾ:2 . ಆರಾ	•	Thu Feb	25 14:58:	(B 1993 _.	10	
573	57 CLN	С	1.3621	3 -17 94055	3 113.59025	λ1 57	
574	57 GLN	_	-40.1897	1 -17.68053	113.34773	A1 57	
575	58 GLY	ัท	-41.8882	5 -19.15742	2 113.41319	8د د۸	
576	58 GLY		-42.8477	1 -19.33514	113.64847	λ1 58	
577	58 GLY		-41.0537	9 -20.23393	3 112.87374	A1 58 A1 58	
578	58 GLY		-40.45534	4 -19.92/01 4 -20 01021	111.50957	A1 58	· · · · · · · · · · · · · · · · · · ·
579 580	58 GLY 59 ALA	о и	-41 34654	4 -19.52227	110.60018	λ1 59	
581	59 ALA		-42.32072	-19.46762	110.83485		0.00000
582	59 ALA	CA	-40.88666	-19.14537	109.26181	A1 59	0.00000
583	59 ALA	CB	-42,07124	-18.74086	108.38101	A1 59	
584	59 ALA	С	-39.86090	-18.02257	109.26073	A1 59	
585	59 ALA	Ο,	-38.85187	-18.04905	108.56603	A1 59	
586	60 LEU	N	-40.12979	-17.02925	110.11113	A1 60	
587	60 LEU	H	-40.96678	-17.03715	110.66448	A1 60	0.00000
588	60 LEU	CA	-39.17026	, -15.9 <i>3</i> 0/9	110.21454	A1 60	0.00000
589 590	60 LEU	CB CG	-41 00342	-14.72100	110.06267	A1 60	0.0000
591	60 LEU	CD1	-41.72077	-13.05637	110.83175	A1 60	0.0000
592	60 LEU	CD2	-40.54968	-13:69403	108.67852	Al 60	0.0000
593	60 LEU	С	-37.86300	16.28607	110.90549	A1 60	0.00000
594	60 LEU	0	-36.81366	-15.71151	110.64266	A1 60	0.00000
595	61 ALA	N	-37.92548	17.30628	. 111.76650	A1 61	0.00000
596	61 AIA	H .	-38.80416	-17.70206	112.04737	A1 61 A1 61	0.00000
597	61 ALA	CA	~36.66060	10 071/7	112.25036 113.37402		0.00000
59E 599	61 ALA 61 ALA	CB C	-35 86652	-18 52303	111.13575	A1 61	0.00000
600	61 ALA	. 0	-34.67753	-18.28483	110.93578	A1 61	0.00000
601	62 A9N	N	-36.59182	-19.33811	110.35468	A1 62	0.0000
602	62 ASN	ĸ	-37.55651	-19.52444	110.56458	A1 62	0.00000
603	62 ASN	CA .	-35.93048	-19.97053	109.20954	A1 62	0.00000
604	62 ASN	ĊВ	-36.90608	-20.83397	108.41185	A1 62	0.00000
605	62 ASN	CG	-36.14296	-21.97501	107.76767	A1 62	0.00000 0.00000
606	62 ASN	OD1	-35.90083	-23.00051	108.38915	Al 62 Al 62	0.00000
607	62 ASN	ND2	-35.81296	-21.80383	106.49294		0.00000
608 609	62 ASN 62 ASN	HD21 HD22	-35.51013	-22.62041	105.92778	2.1 62	0.00000
610	62 ASN	c	-35.27272	-18.97317	108.27635	A1 62	0.00000
611	62 ASN	ō	-34.08977	-19.05772	107.98073	ሕ 1 62	0.0000
612	63 ILE	N	-36.07385	-17.96130	107.91224	A1 63	0.00000
613	63 ILE	H	-37.03805	-17.97906	108.18770	kl 63	0.00000
614	63 ILE	CA	-35.60960	-16.86395	107.05550	A1 63	0.00000
615.	63 ILE	CB	-36.79680	-15.88630	106.84927	A1 63	0.00000
616 617	63 ILE	CG2	-35.71800	-15.60068	105.36486		0.00000
618	63 ILE	CD	-36.27181	-14.79756	105.06416	X1 63	0.00000
619	63 ILE	Ċ	-34.32421	-16.14412	107.48562	k1 63	0.00000
620	63 ILE	0	-33.67028	-15.43835	106.72047	k1 63	0.00000
621	64 ALA	H	-33.97867	-16.34078	108.76481	11 64	0.00000 0.00000
622	64 ALA	Я	=34.55914	-16.88095	109.37800	kl 64 kl 64	0.00000
623	אָדַאָּ 64	CA	-32.68252	-15.86370	109.23001		0.00000
624	64 ALA	CB	+32.78414	-15.37232	110.67448		0.00000
625 626	64 ALA 64 ALA	C 0	-34.53324	-16.51330	108.75412	A1 64	0.0000
627	65 VAL	N	-31 96250	-18.15367	109.50633	A1 65	0.0000
628	65 VAL	н	-32.92093	-18.37377	109.71401	k1 65	0.0000
625	65 VAL	CA.	-30.91507	-19.17596	109.47932	h1 65	0.00000
630	65 VAL	СЭ	-31.28412	-20.41152	110.33210	11 65	0.00000
631	65 VAL	CG1	-31.53866	-19.97521	111.77524 /	41 65	0.00000
632	65 VAL	CG2	-32.45603	-21.23819	109.79675	41 65 41 65	0.00000 0.00000
633	65 VAL	C	-30.45713	-19.56758	108.06237 7		0.00000
634	65 VAL	0	-24.26566	-19 KE037	107.15605	42 66	0.0000
635 636	66 ASP 66 ASP	K E	-32.30553	-19.49723	107.30932	.1 66	0.0000
·	JU MOF	••					

FIG. 11

PCT/US94/05697

_شمدار.	تنت. 2، تنت		جدد لاحق	25 14:58:4	.8 195 <u>3</u>		13	
63,7	66 ASP	CX	11.04858	-19.93216	105.76795	λl	76 6	000000
638	66 ASP	CB			104.93007		66	0.0000
537	SE ASP	CG			105.08208		6 6	0.0000
640	66 ASP	OD1	-33.16986	5 -18.05114	105.20729	Y3	66	0.00000
641	66 ASP	OD2			105.08126		66	0.00000
642	66 ASP	С	-30.18178	-18.83449	105.15564	Y3	66	0.00000
643	66 ASP	0	-29.16482	2 -19.06872	104.50705	YJ	66	0.00000
644	67 LYS	N			105.48017		67	0.00000
645	67 LYS	H		-17.48655			67	0.00000
646	67 LYS	CA	-29.72959	-16.43204	105.19320		67	0.00000
647	67 LYS	CB			105.87104		67	0.00000
648	67 LYS	CG	-29.79159	-13.86177	105.56853		67	0.00000
649	67 LYS	CD	-30.51506	-12.79056	106.38164	XI	67	0.00000
650	67 LYS	CE		-11.38645			67 67	0.00000
651	67 LYS	NZ		-10.44891		V.	67	0.00000
652	67 LYS	HZ1	-30.33262	-9.48398	106.83909		67	0.00000
653	67 LYS	HZ2	-31.70143	-10.49161	108.01156	2)	67	0.00000
654	67 LYS	H23	-30.33221	-16 SADO3	105.64383	Al	67	0.00000
655 656	67 LYS 67 LYS	. C	27.33559	-16:4407B	104.87661	Αì	67	0.00000
657	68 ALA	N	-28 12520	-16.92373	106.92795	Al	68	0.00000
658	68 ALA	H	-28.92119	-17.00935	107.53692	X 1	68	0.00000
659	68 ALA	C)	-26.76352	-17.18143	107.40958	λl	68	0.00000
660	68 ALA	CB	-26.77377	-17.44846	108.91534	A1	68	0.00000
661	68 ALA	C	-26.07149	-18.34364	106.70618	A :	68	0.00000
662	68 ALA	ō.	-24.80989	-18.33297	106.37143	Al	68	0.00000
663	69 ASN	N	-26.87877	-19.37475	106.44973	٨l	69	0.00000
664	69 ASN	н	-27.84416	-19.34421	106.72158	λì	69	0.00000
665	69 ASN	CA	-26.32826	-20.54731	105.77098	ΥJ	69	0.00000
666	69 ASN	CB	-27.33794	-21.70567	105.74618	λì	69	0.0000
667	69 ASN	CG	-27.75534	-22.20215	107.12937	A1	69	0.00000
668	69 ASN	OD1	-28.81753	-22.77967	107.30600	Al	69	0.00000
669	69 ASN	ND2	-26.90880	-21.98927	108.13718		69	0.00000
670	69 ASN	HD21	-26.02949	-21.53117	108.03476	A2	69	0.00000
671	69 ASN	HD22	-27.17968	-22.29754	109.04652	Al	69	0.00000
672	69 ASN	С	-25.83413	-20.26827	104.36379	A1	69	0.00000
673	69 ASN	0	-24.88019	-20.87816	103.89106	A1	69 70	0.00000
674	70 LEU		-26.46696	-19.27268	103.71664	Y.	70	0.00000
675	70 LEU	H	-27.27121	-18.83110	104.12686	71	70	0.0000
676	70 LEU	CA	-25.93555 -26.70466	-10.00313		λl	70	0.00000
677 678	70 LEU	CB CG	-28.07464				70	0.00000
679	70 LEU	CD1	-28 90878	-16 60756	101.23109		70	0.00000
680	70 LEU		-27.93286		99.96378	Al	70	0.00000
681	70 LEU	C	-24.47328	-18.42736			70	0.00000
682	70 LEU	ō	-23.64160	-18.86456	101.72791	Al	70	0.00000
683	71 GLU	N	-24.17065	-17.62592	103.54240	A1	71	0.00000
684	71 GLU	н	-24.87529	-17.33320	104.19342	Al	71	0.00000
685	71 GLU	CA.	-22.77384	-17.24207	103.73624	λl	71	0.00000
686	71 GLU	CB	-22.68099	-16.23884	104.88750	Al	71	0.00000
687	71 GLU	CG	-21.33647	-15.50613	104.93364	YJ	71	0.00000
688	71 GLU	CD	-21.30052	-14.53829	106.10023	A1	71.	0.00000
689	71 GLU	0E1	-20.27115		106.76970	۲ĭ	71	0.00000
690	71 GLU	0E2	-22.29376		106.33423		71	0.00000
691	71 GLU	C,	-21.86369	-18.43808	103.97868	Al	71	0.00000
692	71 GLU	ο,	-20.81243	-18.61050	103.36820	7.1 K.1	71 72	0.00000
693	72 ILE	ĸ	-22.34609	-15.31836	105.86364	7.5		0.00000
694	72 ILE	H	-23.21301	~19.12129			72	0.00000
695	72 ILE	C.F.	-21.56703	-20.33142	105.13649		72	0.00000
€96	72 ILE	CE			106.17923	51	72	0.00000
697	72 115	CG3	-21.51796 -22.55172	-24.00046	107 45409	A.2	72	0.0000
696	72 ILE	CG1	-23.34520	-21 36778	108.51186	Al	72	0.00000
699 730	72 ILE 72 ILE	C.	-21.22106	-21,35113	103.89490	Al	72	0.00000
	72 222	C	21.22300					

FIG. 12

غندر.	ريد . 20شعر	Š	The Year	5 25 14:58:	48 1993		12	
701	72 11	0 3	20.066	70 -21 6816	2 103.64178	Al	72	0.00000
702	73 ME	T N	-22.2541	06 -21.6712	7 103.09738	Al	73	0.00000
763	73 .Mg				2 103.27303		73	0.00000
704	73 ME:		-21.885	77 -22.5114	6 101.95532	λ1	73	0.00000
705	73 ME:	r CB			1 101.56487		73	0.00000
706		r CG		-	2 102.69030		73	0.00000
707	73 ME				2 102.19686		73	0.00000
708	73 ME2			5 -26.5673			73	0.00000
709	73 ME2				0 100.75010		73	0.00000
710	73 ME2			0 -22.3693			73	0.00000
711	74 THE	N	-21.6034	2 -20.4921	4 100.63510	A1	74	. 0.00000
712	74 THE			5 -20.0440			74	0.00000
713	74 THR	C C		3 -19.7383		Al	74	0.00000
. 714	74 THR	CB	-21.4629	9 -18.3549	99.36040	A1	74	0.00000
715	74 THR	. OG1		1 -17.84761		λl	74	0.00000
716	74 THR	. HG1	-20.0469	9 -17.88383	98.02785	λl	74	0.00000
-717	74 TIIR	CG2			100.44667	Al	74	0.0000
718	. 74 THR	. с	-19.3517	0 -19.61359	100.02467	Al	74	0.00000
719	74 THR	0	-18.4855	4 -19.49348	99.16750	λl	74	0.00000
720	75 LYS	·N	·· -19.0853	8 -19.69260	··· 101.33717·	Al	75	0.00000
721	75 LYS	н	-19.8107	5 -19.65220	102.02321	λl	75	0.00000
722	75 LYS	CX	-17.6983	3 -19.83955	101.77078	Al	75	0.00000
.723	75 LYS	CB	-17. 6140	B19.57287	103.27797	λl	75	0.0000
724.	75 LYS	CG			103.86107		75	0.00000
725	75 LYS	CD	-16.2109	1 -19.57437	105.38567	A1	75	0.00000
726	75 LYS	CE	-14.8150	-19.74395	105.98417	Al	75	0.0000
727	75 LYS	NZ			107.44939		75	0.00000
728	75 LYS	HZ1	-13.96820	-19.83508	107.86470	Al	75	0.00000
7.29	75 LYS	HZ2	-15.29159	-18.77551	107.74094	Al	75	0.0000
730	75 LYS	HZ3			107.77639		75	0.00000
731	75 LYS	C.	-17.14118	-21.21778	101.45102	λl	75	0.00000
732	75 LYS	0			100.79660		75	000000
733	76 ARG	N	-17.86151	-22.25466	101.90990	A1	76	0.00000
734	76 ARG	н	-18.69512	-22.09296	102.44436	Al .	76	0.00000
735	7 ARG	CA			101.67525		76	0.00000
736	76 ARG	CB	-18.33654	-24.68494	102.17145	A1	76	0.00000
737	76 ARG	CG	-18.82789	-24.66219	103.62361		76	0.00000
738	76 ARG	CD		-25.98411			76	0.00000
739	76 ARG	NE,		-25.99642			76	0.00000
740	76 ARG	KE		-25.85364		Al	76	0.0000
7.41	76 ARG	CZ			106.34915		76	0.00000
742	76 ARG	NH1		-26.45770			76	0.00000
743	76 ARG	RELL		-26.64144	108.30551		76	0.00000
744	76 ARG	HH12			107.26619		76	0.00000
745	76 ARG	NH2			106.51158		76	0.00000
746 747	76 ARG				107.39854		76 76	0.00000
748	76 ARG				105.72237		76 76	0.00000 0.00000
749	76 ARG	C	-17.00072	-23.91/61	100.20901		76	0.00000
750	77 SER	0			99.83552		77	0.00000
751	77. SER	N	-18.01091		99.36716 2 99.69917 2		77	0.00000
752	77 SER	H, CA	-18.84286		97.93702		77	0.00000
753	77 SER	CB	-17.80506 -19.13837		97.32351		77	0.00000
754	77 SER		-18.92403		96.24728		77	0.00000
755	77 SER	og Hg	-18.99574		95.39572		77	0.00000
756	77 SER	C HG	-17.22854		.97.20137		77	0.00000
757	77 SER	Ö	-17.22834		96.00106		77	0.00000
758	78 ASN	F	-16.52714		97.97646 2		78	0.00000
759	78 ASK	H	-16.44843		98.96113		78	0.00000
76C	78 ASH	CA	-15.74721		97.45264		78	0.00000
761	78 ASN		-14.33390		97.09195		7 E	0.00000
762	76 ASN	CG	-13.52905		98.36114		70	0.02000
763		CSI	-12.89966		98.89264 2		78	0.00000
754		ND2	-13.55264		98.85326		78	0.00000
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./221_	בגם. כאם)	Thu Feb	25 14:58:4	8 1993	13	
7 £5	78 ASN	HD2	1 4.08412	2 -23.19170	98.43678 Al	78	0.00000
766	78 ASN	н н д 2	2 -13.03873	1 -22.65362	99.68493 Al	78	0.00000
. 7 67	78 ASN	C	-16.34134	1 -19.72172	96.31352 A1	78	0.00000
766	78 ASN	0		3 -19.42979		78	0.00000
769	79 TYR			2 -19.32597	96.52354 A1	79	0.00000
770	79 TYR			-19.55771	97.38252 Al	79	0.00000
771	79 TYR			3 -18.48918	95.49858 Al	79 79	0.00000
772	79 TYR 79 TYR			-18.92649 -20.32764	95.21260 Al 94.65360 Al	79	0.00000
773 774	79 TYR			-21.23278	95.25950 A1	79	0.00000
775	79 TYR			-22.54189	94.75217 Al	79	0.00000
776	79 TYR			-20.72112	93.54108 Al	79	0.00000
777	79 TYR			-22.03339	93.03642 Al	79	0.00000
778	79 TYR			-22.93822	93.64752 A1	79	0.00000
779	79 TYR	ОН	-20.07585	-24.23438	93.18578 A1	79	0.00000
780	79 TYR	нн	-19.67682	-24.30670	92.31232 Al	79	0.00000
781	79 TYR	C		-17.01261	95.83719 A1	79	0.00000
782	79 TYR	0		-16.51273	96.76683 Al	79	0.00000
783	80 THR	N		-16.33235	95.00452 A1	80 80	0.00000 0.00000
784 785	80 THR	Н		-16.79509 -14.92660	94.24288 A1 95.20513 A1	80	0.00000
786	80 THR	CB		-14.87711	96.18098 A1	80	0.00000
787	BO THR	OGI		13.52440	96.43537 A1	80	0.00000
788	BO THR	HG1		-13.50159	96.80560 A1	80	0.00000
789	80 THR	CG2		-15.72265	95.72104 A1	80	0.00000
790	80 THR	С	-16.77512	-14.32258	93.83840 Al	80	0.00000
791	80 THR	0		-14.97644	92.99811 Al	80	0.00000
792	81 PRO	N		-13.08096	93.61451 Al	81	0.00000
793	B1 PRO	CD	-18.04687		94.50529 Al	81	0.00000
794	B1 PRO	CA		-12.43774	92.32255 A1	81	0.00000
795	81 PRO	СВ	-18.08102		92.28996 A1	81 81	0.00000 0.00000
796 797	81 PRO 81 PRO	CC	-18.22970 -15.57247		93.74456 X1 92.25040 A1	81	0.00000
798	81 PRO	Ο .	-15.41926		92.16776 A1	81	Ü.00000
799	82 ILE	N.	-14.55883		92.27154 A1	82	0.00000
800	82 ILE	H	-24.69364		92.31920 A1	82	0.00000
801	82 ILE	CA	-13.18946		92.13921 A1	82	0.00000
802	82 ILE	CB	-12.60010		93.53959 Al	82	0.00000
003	82 ILE	CG2	-12.41140		94.38674 A1	82	0.00000
804	B2 ILE	CG1	-11.31152		93.44331 Al	82	0.00000
805	82 ILE	CD	-10.76554		94.80896 A1 91.37376 A1	82 82	0.00000
806 807	82 ILE 82 ILE	C OCT1	-12.35649 -11.36717		90.75062 A1	82	0.00000
808	82 ILE	OCT2	-12.72556		91.38671 A1	82	0.00000
809	83 GLY			-0.31236		1	0.00000
810	83 GLY	HT1	-17.21994	0.44323	94.35235 B1	1	0.00000
811	83 GLY	HT2	-16.86357	-1.12219	94.96444 Bl	1	0.00000
812	83 GLY	HT3	-17.61098	0.01920	95.97150 Bl	1	0.00000
813	83 GLY	CA	• • • • • • • • • • • • • • • • • • • •	0.91116	94.55151 B1	1	0.00000
814	83 GLY	С	-18.62573	-2.38203	94.66351 B1	1	0.00000 0.00000
815 816	83 GLY 84 ASP	0	-17.35786 -19.57260	2.70920	94.84086 B1 94.59303 B1	1 2	0.00000
817	84 ASP	N H	-20.49658	-3.20239 -2.89510	94.35600 B1	2	0.00000
818	84 ASP	CA	-19.43900	-4.63200	94.86181 B1	2	0.00000
819	84 ASP	СВ	-19.44643	-4.83356	96.38475 B1	2	0.00000
820	84 ASP	CG	-18.89301	-6.19619	96.73815 Bl	2	0.00000
821	04 ASP	OD1	-17.69451	-6.29764	96.98109 Bl	2	0.00000
822	84 ASP	OD2	-19.66566	-7.14958	96.75715 B1	2	0.00000
923	84 ASP	С	-20.62766	-5.31072	94.19848 Bl	2	0.00000
824	84 ASP		-21.46903	-4.61697	93.63479 B1	2	0.00000
925	85 THR		-20.67796	-6.64606	9;.24891 B1	3 3	0.00000 0.00000
926 527	85 THR	н	-20.04362	-7.15926	94.84032 B1 93.59154 Ei	3	0.00000
527 328	65 THR 85 THR	C2	-21.75257 -21.58903	-7.39367 -7.31950	93.39134 E1 92.05122 E1	3	0.00000
	UU ARS	<u></u>		-1.51750		-	

. / تعدر	_KIN2.CR	Đ	Thu Feb	25 14:58:4	B 1993	14		
829	* 85 TH	R OG1	22.762	5 -7.82480	91.39964 61	3		0.00000
830	85 TH		-22.7343			3		0.00000
83%	85 TH	R . CG2	-20.3296			3		0.00000
832	85 TH	R C	-21.7429	0 -8.83241	94.09035 B1	3		0.00000
833	85 TH		-20.7645	4 -9.30377		3		0.00000
834	86 AR		-22.8649			4		0.00000
835	B6 AR		-23.6240			4		0.00000
836	86 ARG			0 -10.87552		4		0.00000
837	86 ARC			8 -10.79083		4		0.00000
838	86 ARC			3 -11.65365	- '	4		0.00000
839 840	86 ARC			6 -13.10294		4		0.00000
841	86 ARG			0 -14.03911 3 -14.68836		4		0.00000
842	86 ARG	-		2 -14.05016		4		0.00000
843	86 ARG			1 -14.85617		4		0.00000
844	86 ARG			5 -14.87872	99.26299 Bl	4		0.0000
845	86 ARG	HH12	-20.0519	4 -15.47498	97.69021 B1	4		0.00000
846	36 ARG			7 -13.26636	99.42563 Bl	4		0.0000
847	86 ARG				100.26897 B1	4		0.00000
848	86 ARG				· 99.36831 B1	4		0.00000
849	86 ARG			2 -11.82065	93.57905 B1	4		0.00000
850 851	86 ARG 87 PRO			0 -11.52883	93.17882 B1	4		0.00000
852	87 PRO			0 -12.98532 9 -13.34562	93.27325 B1 93.54376 B1	5 5		0.00000
853	87 PRO			9 -14.02757	92.53087 B1	5		0.00000
854	87 PRO	CB		-15.13066	92.39999 B1	5		0.00000
855	87 PRO			-14.43376	92.52460 B1	5		0.00000
856	87 PRO	С		-14.54496	93.21975 B1	5	•	0.00000
857	87 PRO	0	-25.32441	-14.39544	94.41838 El	5		0.00000
858	88 ARG	N	-25.94344	-15.17061	92.39123 B1	6		0.00000
859	88 ARG	н		-15.35588	91.44433 B1	6		0.00000
860	88 ARG	CA		-15.62675	92.90235 B1	6		0.00000
861	80 ARG	C3		-15.39805	91.81653 B1	6		0.00000
862 863	88 ARG	CD CD		-15.55859 -14.91787	92.35844 B1 91.49236 B1	6 6		0.00000
864	88 ARG	NE NE		-14.57277	92.34607 B1	6 .		0.00000
865	88 ARG	HE .		-14.64038	93.33788 B1	6		0.00000
866	88 ARG	CZ		-14.12165	91.86193 B1	6		0.00000
8 67	88 ARG	йнэ	-34.06565		92.71322 B1	6		0.00000
8 6 B	88 ARG			-13.48197	92.40144 B1	6		0.00000
869	88 ARG			-13.92533	93.69838 B1	6		0.00000
870 871	88 ARG			-13.97383	90.54974 B1	6		0.00000
872	88 ARG 88 ARG			-13.64013 -14.19859	90.16725 B1 89.92214 B1	6 6		0.00000 · 0.00000
873	BB ARG			-17.07404	93.36366 B1	6		0.00000
874	88 ARG	o ·		-18.00090	92.63574 B1	6		0.00000
875	89 PHE			-17.24057	94.62546 B1	7		0.00000
876	89 PHE		•	-16.46797	95.18173 B1	7		0.0000
877	89 PHE	CA	-27.64368	-18.59443	95.17433 B1	7	i	0.00000
	89 PHE		-26.56379	-18.74887	96.25427 Bl	7		0.0000
879	89 PHZ			-18.55805	95.61829 B1	7		0.0000
880	89 PHE		•	-19.48139	94.64858 B1	7		0.00000
8 <i>8</i> 1 882	89 PHE			-17.43023	95.96102 B1	7		0.00000
882 883	244 68 244 68			-19.27079 -17.21976	94.00961 B1 95.32384 B1	7 7		0.00000 0.00000
884	89 PHE			-17.21976	95.32384 B1 94.34976 B1	7		0.00000
885	89 PHE		-28.99632		95.72084 B1	, י		0.00000
306	89 PHE		-29.85551	-18.12429	95.94786 B1	י ר		0.0000
887	90 LED		-29.15750		95.87791 B1	8		0.0000
886	90 LEU	н -	26.38720	-20.92114	95.76376 Bi	8		0.0000
889	90 LEU		-30.48975		96.14113 B1	9		0.0000
890	90 LEU		31.01265		94.64981 B1	8		0.00000
891	90 LEU			-21.16692		3		0.0000 0.0000
992	90 LEU	CD1 -	32.81137	-21.76230	93.19113 21	3		7.00000

. /ರಸು_	KIN2.CAD	•	Thu Feb	25 14:58:4	8 1993 _.	15	
893	. 90 TEA	CD2	3.4749	B -21.40699	95.55222 B	. 8	0.00000
894	90 LEU	C		7 -21.90281		. 8	0.00000
دوة	90 LEU	0	29.7421	6 -22.89200	97.07224 BI		0.00000
896	91 TRP	N	-31.2813	3 -21.71086			0.00000
897	91 TRP	ĸ		7 -20.85226			0.00000
898	91 TRP	CY		7 -22.06043			0.00000
899	91 TRP	CB	-31.76159	9 -22.50789	100.54418 B1	. 9	0.00000
900	91 TRP	cc	-30.46050	-22.11490	101.18157 B1	. 9 . 9	0.00000
901	91 TRP	CD2	-29.22893	-22.79283	101.16170 B		0.00000
902		CE2			101.98847 B1		0.00000
903	91 TRP	CE3		-24.00327 -20.96427			0.00000
.904	91 TRP 91 TRP	CD1	-30.26500) -20.96427 \ -20.96427	102.44030 B1		0.00000
905 906	91 TRP 91 TRP	HE1		-20.15616			0.00000
907	91 TRP	C22		-22.40520			0.00000
908	91 TRP	CZ3			100.81577 B1		0.00000
909	91 TRP	CH2	-26.56124	-23.61859	101.58828 B1	9 .	0.00000
910	91 TRP		-32.96765	-23.39534	98.70806 B1	9	0.00000
911	91 TRP	0		-22.66026	98.49798 Bl	9	0.00000
912	92 GLN -	-N	33.02847	-24.71550		10	0.00000
913	92 GLN	H	-32.20362	•		10	0.00000
914	92 GLN	CA		-25.34751		10	0.00000
915	92 GLN	CB		,-,26.04119		10	0.00000
916	92 GLN	CG		-25.05724		10	0.00000
917	92 GLN	CD		-25.79390		10	0.00000 0.00000
918	92 GLN	OE1	-32.99424			10 10	0.00000
919	92 GIN	NE2		-26.32196		10	0.00000
920	92 GLN	HE21		-26.21874	94.65013 B1 93.40658 B1	10	0.00000
921 922	92.GLN	HE22		-26.83019 -26.35610	99.77327 B1	10	0.00000
922	92 GLN 92 GLN	C 0		-27.07419		10	0.00000
924	93 LEU	N .	-35.49484	-26.39205		11	0.00000
925	93 LEU	Н				11	0.00000
926	93 LEU	Cγ	-35.59559	-27.43909	101.48633 B1	11	0.00000
927	93 LEU	CB			102.88181 B1	11	0.00000
928	93 LEU	CG.		-27.70766	103.99848 Bl	11	0.00000
929	93 LEU	CD1	-34.33031		105.14843 Bl	11	0.00000
930		CD2	-35.82812	-28.74167	104.52805 B1	11	0.00000
931		C	-36.93596	-28.11112	101.32695 B1	11	0.00000 0.00000
932		0		-27.47571	100.96621 Bi 101.56549 Bl	11 12	0.00000
933		N.	-36.95045		101.56549 B1 101.79736 B1	12	0.00000
934		H		-29.89066 -30.14121		12	0.00000
935 936	•	CA CB	-30.21003	-30.14121	100.13839 B1	12	0.00000
937		CC	-30 43761	-31.85442	100.07544 B1	12	0.00000
938		æ	-39.91969	-32.44832	98.75113 H1	12	0.00000
939		CE	-41.00198		99.23924 B1	12	0.00000
940		NZ	-41.98389	-33.90963	98.26846 Bl	12	0.00000
941	94 LYS	K21			98.86351 B1	12	0.00000
942		HŻ2	-42.49249	-33.14263	97.78947 B1	12	0.00000
943		HZ3	-41.57162	-34.57937	97.59540 Bl	12	0.00000
944		С	-38.34910	-31.08699	102.67642 Bl	12	0.00000
945		0	-37.54720	-31.99770	102.85849 B1	12 13	0.00000
946	•	N.	-39.40676	-30.85929	103.46609 Bl	13	0.00000
947 948	•	H	-39.99321	-30.05201	103.34519 Bl 104.49053 Bl	13	0.00000
949		CA CB	-37./3/34	-31 40224	105.86481 E1	13	0.00000
950	•	CG	-38.13531	-30 20819	106.49960 B1	13	0.00000
951		CD1	-39 84517	-28.96286	106.12996 B1	13	0.00000
952		CD2	-40.60668	-30.26479	107.51173 B1	13	0.00000
953	•	CE 1	-39.45587	-27.78326	106.77985 B1	13	0.00000
954		CE2	-41.01477	-29.08019	108.16313 B1	13	0.00000
955	95 PHE 0	22	-40.43557	-27.84425	107.79605 21	13	0.00000
956	95 PME (3	-41.22005	-32.16799	104.57423 E1	13	0.00000

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95. 750	95 P	HE O	- 08591	-31,31332	104.41691	B1 1	3 -0::00000
958	96 G		47812	-33.46734	104.75822	B1 1	0.00000
9: 9	76 G		-40.74093	-34.09871	105.01948	B1 1	
960	96 G		-42.80002	-33.99059	104.40078	B1 1	
961	96 G	LU CB	-42.75085	-34.25375	102.88321	B1 1	
962	96 G		-43.92070	-34.87965	102.11/91	B1 1 B1 1	•
963	96 G		-43.55505	-34.92496 -34.15253	99.84297	B1 1	
964	96 G		-44.08203	-35.68559	100.23154	B1 1	
965 966	96 G		-43.13129		105.17925	B1 1	4 0.00000
967	96 G		-42 27510	-36.11395	105.37879	B1 1	
968	97 C		-44 39621	-35.34431	105.62501	B1 1	
969	97 C		-45.05281	-34.60566	105.44206	B1 1	
970	97 C	YS CA	-44.78990	-36.60747	106,25700	B1 · 1	
971	97 C	YS CB	-45.09404	-36.46870	107.76554	B1 1	
972	97 C		-46.49873	-35.54248	108.42288	B1 1	- · · · · · · · · · · · · · · · · · · ·
973	97 C		-45.89248	-37.36966	105.55700	B1 1	_
974	97 C		-46.75253	-36.80546 -38.69951	104.00707		-
975	98 H.		-45.8U350	-39.11332	105.71704	_	
976	98-H		-45.06634	-39.59530	105.05785		
977 978	98 HI 98 HI		-46 00867	-40.61029	104.19437	B1 1	6 0.00000
979	98 H		-45.38253	39.98357	102.97364	B1 1	6 0.00000
980	98 H		-45 88967	-40.13539	101.74183	B1 1	
981	98 HI		-46.72307	-40.60043	101.51181	B1 1	
982	98 H		-44.21286	~39.22166	102.89430	B1 1	
983	98 H	S NE2	-44.02278	-38.92334	101.58753	B1 1	
984	98 KI		-45.0526B	-39.48082	100.87633	B1 1 B1 1	•
985	98 KI		-47.56750	-40.40062	106.05217		•
986	98 H		-47.04279	-40.99638 -40.39547	105.33223		=
987	59 51				105.00666	B1 1	
988	99 PI		-49.25243	-41.01889	106.77662		
989 990	99 PI		-50.48946		107.59234	B1 1	7 0.00000
991	99 P.		-49.63868	-39.72312	108.81497	B1 1	0.00000
992	99 PH		-48.50685	-38.87580	108.76543	B1 1	7 0.00000
993	99 PH		-49.92563	-40.47457	109.97439	B1 1	
994	99 PH		-47.64103	-38.80002	109.87537	B1 1	
995	99 PH	E CE2	-49.06252	-40.39779	111.08354	B1 1	
996	99 PH		-47.92137	-39.56658	111.02626	B1 1	
997	99 PH			-41.97404 -41.71068	105.19800		
998	99 PI	•	-51.48007 -50.86837		106 85844		
999 1000	100 PH		-50.32768	=43.27017	107.68944	B1 1	
1001	100 PH	•	-51.84718	-44.10783	106.41132	B1 1	_
1002	100 PH	•	-51.44468	-45.52210	106.84262	B1 1	
1003	100 PH		-51.08740	-46.34931	105.63001	BI I	
1004	100 PH	E CD1	-49.81159	-46.95065	105.55366	B1 1	
1005	100 PH		-52.01269	-46.50158	104.56911	B1 1	-
1006	100 PH		-49.45012	-47.69723	104.41336		
1007	100 PH		-51.65494	-47.24829 -47.83859	103.42413	B1 1	
•	100 PH		-50.3/400	-43.84401	106.96975	B1 1	
1009 1010	100 PH		-53 38235	-43.49692	108.13319	B1 1	
1011	100 FA		-54 19611	-44.08275	106.08672	91 1	
1012	101 AS		-53.92777	-44.28429	105.14366	B1 1	
1013	101 AS		-55.63451	-43.97453	106.37273	B1 1	
1014	101 አ5		-56.35400	-45,03359	105.52094	P1 1	
1015	101 AS		-57.86040	-44.83624	105.52911	21 1	
1016	101 AS		-58.43246	-44.06601	104.77523	21 1	
1017	101 AS		-58.51327	-45.52184	100.41333		•
1018	101 AS		-58.04843	-46.21753 -45.49785	106.46882		
1015	101 AS		-39.30000	-45.49763	107.83398	B1 1	
1020	101 45	:: c	-39.00277	-, -, -, -, -, -, -, -, -, -, -, -, -, -			

./521	تبتته"	.೮೮		Thu Feb	25 14:58:	4E 1993		17	
1021	101	ASN	٥	: 6 7373	7 -43.2399	7 108.39463	B1	19	0.00000
1022		GLY	ĸ	- 25.6346				20	0.00000
1005	202	GLY	H	-55.0037	0 -45.8218			20	0.00000
1024	102	CLY	CA	-55.9743	8 -45.3235			20	0.00000
1025		CLY	С	-55.0208				20	0.00000
1026	102	GLY	0	-54.4600		4 111.73800		20	0.00000
1027	103	THR	N	-54.8474		6 110.54648		21	0.00000
1028	103	THR	Н			1 109.77435		21	0.00000
1029	103	THR	CA			2 111.30363		21 21	0.00000
1030	103	THR	CB	-54.7548				21	0.00000
1031 1032	103 103	THR THR	0G1	-54.8197	5 -40.5335 4 -40.0991			21	0.00000
1032	103	THR	HG1 CG2		5 -42.7243	-		21	0.00000
1034		THR	C		3 -42.9238		Bl	21	0.00000
1035		THR	Ö		6 -42.5928			21	0.00000
1036		GLU	N		7 -43.7843	3 110.78327	Bl	22	0.00000
1037	104	GLU	11	-52.5083	-43.9402	1 109.90068	Bl	22	0.00000
1038	104		CA		-44.4245			22	0.00000
1039	104		CB		7 -45.9519			22	0.00000
1040						5 112.41591		22	0.00000
1041	104		CD			113.43075		22 22	0.00000
1042	104		0E1 0E2			113.08359		22	0.00000
1044	104		C		-43.91901			22	0.00000
1045		GLU	0	-49.58628			Bl	22	0.00000
1046		ARĢ	N		-44.04578		Bl	23	0.00000
1047		ARG	H	-48.46820				23	0.00000
1048		ARG	CA	-47.23862				23	0.00000
1049		ARG	CS		-43.14957		Bl	23	0.00000
1050	105	AP.Ġ	CG	-45.11167			Bl	23	G.00000
1051	105	ARG	CD	-43.89570	-43.22016			23	0.00000
1052		ARG	ΝE	-42.65640				23	0.00000
1053		ARG	HΞ			112.83941		23	0.00000 0.00000
1054	105		CZ			110.93892		23 23	0.00000
1055 1056	105 3		221 221	-41.02352	-41.00938 -40.49136			23	0.00000
1057	105 3		HH12		-40.92747			23	0.00000
1058	105 2		NH2		-41.82718			23	0.00000
1059	105 7		HH21		-41.15340			23	0.00000
1060	105 7	LRG	RH22			109.38941	Bl	23	0.00000
1061	105 3		Ċ				B1	23	0.00000
1062	105 }		0				Bl	23	0.00000
1063	•		N 		-43.51517		Bl Bl	24 24	0.00000
1064	106 V		H Ch	-46.72693	-42.61352	107.58352		24	0.00000
1066	106 V		CA CB	-46 27081	-44.39073	106.24298	B1	24	0.00000
1067	106 V			-45.79579	-45.65575	105.52366	B1	24	0.00000
1068	106 V		CG2	-47.77990	-44.41166	106.42922	Bl	24	0.00000
1069	106 V	'nĹ (C	-44.14065	-43.66075	107.41554	B1	24	0.00000
1070	• • •	•	0			108.29696		24	0.00000
1071	107 A		N	-43.87314	-42.98069	106.29416	B1	25	0.00000
1072	107 A		H	-44.58452	-42.69040	105.65671	B1	25 .	0.00000
1073	107 አ		CA			106.12329		25	0.00000 0.00000
1074	107 A		CB	-41.95685	-43.01317	104.76032	21	25 25	0.00000
1075 1076	107 A		CG CD	-41.96328	-42.02619	103.58362	B)	25 25	0.00000
1077	107 A		ZD ZD	-42.3333/ -43 76939	-42.00029 -27 C3515	102.23147	B1	25	0.00000
1078	107 A		IE	-44.37946	-42.14312	102.28284	Bl	25	0.00000
1079	107 A	•	22	-41.26510	-44.17490	102.20945	Bl	25	0.0000
1080	107 A	RG :	VH 1	-45.56598	-44.32747	102.35905	E :	25	0.0000
1081	107 X	RG :	:H11	-46.08629	-45.19207	102.34562	91	25	0.00300
1082	107 A		1112	-45.19539	-43.55730	102.54612	E 3	25	0.00000
1093	107 A	RG :	:::2 	-43.47983	-45.23695	102.05798	21 21	25 25	0.00000 0.00000
1084	107 A	?.G :	HE21	-43.86541	-45.15896	102.05452	= 1		0.0000

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1085	107 ARG	нн22	2 49404	-45.11997	101.94496	Bl	25	0.00000
1086	107 ARG	C	2 37787	-41.03546	106.32504	Bl	25	0.00000
1007	157 ARG	ō	-43.36896	-40.31981	106.43645	Bl	25	0.00000
1088	108 LEU	N	-41,12633	-40.58602	106.39627	Bl	26	0.00000
1089	108 LEU	H	- 40 - 32617	-41.17985	106.29320	Bl	26	0.00000
1090	108 LEU	CA	-40.89942	-39.15744	106.53751	81	26 26	0.00000
1091	108 TEU	CB	-40.80087	-38.80954	108.01248 108.35874	B1	26	0.00000
1092	108 LEU	CG	-41.9634/	-38 AA720	109.53808	B1	26	0.00000
1093	108 LEU	CD2	-41 50070	-36 44650	108.50151	Bl	26	0.00000
1095	108 LEU	C	-39 62061	-38.77273	105.85764	Bl	26	0.00000
1095	108 LEU	ō	-38.76368	-39.61093	105.60154	B1	26	0.0000.0
1097	109 LEU	N	-39.52150	-37.47828	105.56773	Bl	27	0.00000
1098	109 LEU	н	-40.28459	-36.84167	105.72467	B1	27	0.00000
1099	109 LEU	CA	-38,30683	-37.02679	104.90334	Bl	27	0.00000
1100	109 LEU	CB	-38.47418	-37.24039	103.39009	BI	27 27	0.00000 0.00000
1101	109 LEU	CG	-39.74490	-36.59695	102.82634	B.7	27	0.00000
1102	109 LEU	CD1	-39.42210	-35.25631	102.17066	BI	27	0.00000
1103	109 LEU 109-LEU	CD2	-40.49190	-37.50323	-105.21268	Bl	27	0.00000
1104 1105	109-LEU	0	-34 AR411	-34.77322	105.47943	Bl	27	0.00000
1105	110 GLU	ห	-36.70092	-35.29768	105.12126	Bl	20	0.00000
1107	110 GLU	11	-36.01498	-36.01786	105.00992	Bl	28	0.0000
1108	110 GLU	CV	-36.20315	-33.92985	105.08184	Bl	28	0.00000
1109	110 GLU	CB	-35,80977	-33.47627	106.49578	Bl	28	0.0000
1110	110 GLU	CĢ	-35.32216	-32.02210	106.60993	Bl	28	0.00000
1111	110 GLU	CD	-33.83808	-31.87178	106.30413	Bl	28	0.00000
1112	110 GLÚ	OEl	-33.37957	-30.74951	106.10521	Bl	28	0.00000
1113	110 GLU	OE2	-33.09671	-32.85201	106.36491	Bl	28	0.00000
1114	110 GLU	C	-35.01026	-33.94076	104.15500	B1	28 28	0.00000
1115	110 GLU	0	-34.25577	-34.90807	104.11585	בם בת	29	0.00000
1116	111 ARG	N	-34.89439	-32.86843	103.37222	B1	29	0.00000
1117	111 ARG	H	-35.61365 -33.75665	-32.16766	102.46249	B1	29	0.00000
1118	111 ARG 111 ARG	CA CB	-33.99004	-33.60407	101.20352	B1	29	0.00000
1119 1120	111 ARG	ĊG	-35.45106		100.75305	Bl	29	0.0000
1121	111 ARG	CD	-35.67880		99.75174	Bl	29	0.00000
1122	111 ARG	NE	-37.09124	-35.21147	99.72614	B1	29	0.0000
1123	111 ARG	HE	-37.71860		100.22757		29	0.00000
1124	111 ARG	CZ	-37.50098	-36.32384	99.09789		29 29	0.00000
1125	111 ARG	NH1	-38.78116	-36.68971	99.16119 98.68276		29	0.00000
1126	111 ARG	нн11	-39.12269	-37.49888	99.70940		29	0.00000
1127	111 ARG		-39.43712 -36.63391	-30.16093	98.41518	_	29	0.00000
1128	111 ARG	NH2	-36.91501	-37 90348			29	0.00000
1129 1130	111 ARG	HH22	-35.67503	-36.78688	.98.36264	B1	29	0.00000
1131	111 ARG	С	-33.49135	-31.31194	102.08170	Bl	29	0.0000
1132	111 ARG	ō	-34.39395	-30.54569	101.76699	31	29	0.00000
1133	112 CYS	N	-32,20568	-30.97057	102.12414	Bl	30	0.00000
1134	112 CYS	H .	-31.50733	-31.64380	102.37455	Bl	30	0.00000
1135	112 CYS	CA	-31.80468	-29.62360	101.73826	B1	30	0.00000 0.00000
1136	112 CYS	CB	-31.12874	-28.92365	102.91930	B.1	30 30	0.00000
1137	112 CYS	5G	-30.70297	-27.19555	100.54300	B)	30	0.00000
1138	112 CYS	С 0	-30.87388 -29.97769				30	0.00000
11,39 1140	112 CYS 113 ILE	N .	-31.15975	-28.70696	99.65078	B1	31	0.00000
1141	113 ILE	H :	-31.84193	-28.00116	99.86111	51	31	0.00000
1142	113 ILE	CA.	-30.55306	-28.70228	98.32464	Bl	31	0.00000
1143	113 ILE	C3	-31.56021	-29.33526	97.30706	Bl	31	0.00000
1144	113 ILE	CG2	-33.00339	-29.37434	97.82323		31	0.00000
1145	113 ILE	CG1	-31.52431	-28.74184	95.89635		31	0.00000
1146	113 ILE	CD	-32.44047	-29.50025	94.93272		31 31	0.00000 0.00000
1147	113 ILE	C	-30.08576	-27.29665	97.96536 98.21317		31	0.00000
::48	113 ILE	C	-30.75333	-25.25900	yo. 2131/	<i>.</i>	J 1	0.0000

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1149	11	4 TYR	и	28.8701	0 -27.2556	4 97.41665 B	32	0.00000
1150	` 11		-		0 -28.1015	1 97.10297 B	32	0.00000
1151	11	4 IYR	CA	-28.2203	6 -25.9826	4 97.11462 B	1 32	0.00000
1152	11				8 -26.0193		32	0.00000
. 1153	11				7 -24.7210	7 97.61311 B	32	0.00000
1154	11				0 -23.4582			0.00000
1155		4 TYR			6 -22.2808			0.00000
1156	11				4 -24.8021			0.00000
1157 1158		4 TYR 4 TYR			1 -23.6265 D -22.3694			0.00000
1159	11.			•	7 -21.2191			0.00000
1160	11				6 -21.4301			0.00000
1161		TYR	C		-25.8321			0.00000
1162	114		ō		-26.7096			0.00000
1163		ASN	N		-24.6999		. 33	0.00000
1164		ASN	H	-29.03092	-24.02393			0.00000
1165		NZA (CA		-24.41545			0.00000
1166		ASN	CB		-23.84385			0.00000
1167		ASN	CG		-22.43118			0.00000
1168						194.85370 B1 9 92.96271 B1		0.00000 0.00000
1169 1170		NZA i	ND2	-28.04883	-21.48359		_	0.00000
1171		ASN .	HD21					0.00000
1172		ASN	C		-25.59300			0.00000
1173	115		Ö		-25.91402			0.00000
1174		GLN	N		-26,23583	1		0.00000
1175		GLN	н		-25.92614			0.00000
1176		GLN	CA	-30.60575	-27.38897	92.31212 B1	34	0.00000
1177	116	GLŅ	CB		-26.98635			0.0000
1178		GLN	CG		-27.96419		_	0.00000
1179		CTN	CD		-27.31473		34	0.00000
1180	116		OE1		-27.33002			0.00000 0.00000
1181	116		NE2 HE21	-29.39553	-26.71450 -26.71396		34	0.00000
1182 1183		GLN	HE22	:	-26.24641		34	0.00000
1184		GLN	C.		-28.73663		34	0.00000
1185		GLN	0		-29.77893		34	0.00000
1186	117	GLU	N	-28.76921	-28.72803	93.24838 B1	35	0.00000
1187	117	GLU	H	-28.34990	-27.88576		35	0.00000
1188	117	GLU	CA	-28.17324	-30.02538	93.58636 B1	35	0.0000
1139	117	GLŲ	CB		-30.02385	93.23572 Bl	35	0.00000
1190	117		CG	-26.41125	-29.71932	91.75724 B1	35 35	0.00000
1191 1192	117	GLU GLU	CD OE1	-24.93459 -24.62388	-29.87167	91.43636 B1 90.37359 B1	35	0.00000
1193	_	GLU	OE2	-24.09945		92.24135 Bl	35	0.00000
1194		GLÜ		-28.34342	-30.39133		35	0.00000
1195	117	GLU	ο . ΄	-28.39032	-29.54087	95.93593 Bl	35	0.00000
1196		ern	N	-28.45418		95.31151 B1	36	0.0000
1197	118		H	-28.37884		94.57056 B1	36	0.00000
1198	118		CA	-28.64640		96.70504 H1 96.80220 B1	36 36	0.00000 0.00000
1199 1200	118 118	GT. GT.	CB CG	-28.86529 -30.04821		96.80220 B1 96.04454 B1	36	0.00000
1201	118		CD	-30.04621		96.39969 B1	36 ·	
1202	118	GLU	OE1	-31.31366		96.62004 Bl	36	0.00000
1203		ern	022	-29.17775		96.46236 B1	36	0.0000
1204		ern ern	C.	-27.45968		97.59954 B1	36	0.00000
1205	118		0	-26.30375	-31.93771	97.22004 B1	36	0.0000
1206	119		R .	-27.77719	-31.35385	98.81671 B1	37	0.00000
1207	119	SER		-28.73032		99.10568 51	37	0.00000
1208	119		CA	-26.67523	-31.09310	99.74333 B1	37	0.00000
1209	119					100.25438 B1	37 37	0.00000 0.00000
1210	119		03	-25.62406	-29.23340	100.96554 B1 101.30354 51	37	0.00000
1211	119			-25.73940			37	0.00005
	1:7		_	יסיינים. ט ג -	~~	190.0:0:0:	- '	

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1717	- - 119 SER	٥	15 72568	-32.90554	101.03211	Bl	37	0.00000
1213	120 VAL	. ม	-27 72434	-32.06808	101.67561	81	38	0.00000
1215	120 VAL	់អ	-28 50026	-31.45938	101.48344	Bl	38	0.00000
1216	120 VAL	شم	-27.82678	-32.99802	102.80099	Bl	38	0.00000
1217	120 VAL	CB	27.44790	-32.32244	104.13888	Bl	38	0.00000
1218	120 VAL	CG1	-25.93252	-32.18078	104.29388	Bl	38	0.00000
1219	120 VAL	CG2	-28.15631	-30.97900	104.33500	Bl	38	0.00000
1220	120 VAL	С	-29.23777	-33.53625	102.89372	B1	38	0.00000
1221	120 VAL	0 .	-30.19812	-32.88656	102.49711	Bl	38	0.00000
1222	121 ARG	N -5	-29.34164	-34.75356	103.41944	Bl	39 39	0.00000
1223	121 ARG	H	-28.52525	-35.27675	103.67737	D1	39	0.00000
1224	121 ARG	CA	-30.65941	-35.37282	103.49447	D)	39	0.00000
1225	121 ARG	CB	-30.83879	-36.26556	102.25451	B1	39	0.00000
1226	121 ARG	CG	-32.24891	-36.83907	102.13977	ום	39	0.00000
1227	121 ARG	CD	-32.36448	-38.11228 -38.89614	101.31436	BI	39	0.00000
1228	121 ARG	NE	-33.46636	-38.82444	102.86363	Bl	39	0.00000
1229	121 ARG	HE	34.21779	-30.02377	101.13259	Bl	39	0.00000
1230	121 ARG 121 ARG	CZ NH1	-35 18842	-40.40193	101.73316	B 1	39	0.00000
1231	121 ARG	MUT	-35.77971	-41.03741	101.23780	B 1	39	0.00000
1232 1233	121 ARG	нн12	-35.33365	-40 28453	102.71808	Bl	39	0.00000
1233	121 ARG	NH2	-33.99786	-39.84536	99.82444	Bl	39	0.00000
1235	121 ARG		-34.53079		99.25675		39	0.00000
1236	121 ARG	HH22	-33.27054	-39.30635	99.39731		39	0.00000
1237	121 ARG	С	-30.80743	-36.21013	104.75602	Bl	39	0.0000
1238	121 ARG	Ö	-29.86515	-36.82872	105.23483	Bl	39	0.00000
1239	122 PHE	N	-32.0407S	-36.24095	105.26404	Bl	40	0.00000
1240	122 PHE	н	-32.75061	-35.63B10	104.89111	B 1	40	0.00000
1241	122 PHE	CA	-32.40668	-37.26064	106.24769	Bl	40	0.00000
1242	122 PHE	CB	-33.75724	-36.87165	106.85296		40	0.00000 0.00000
1243	122 PHE	CG	-33,64992	-36.46831	108.30418	Bl	40	0.00000
1244	122 PHE	CD1	-32.56100	-35.69712	108.77664	B1	40 40	0.00000
1245	122 PHE	CD2	-34.66985	-36.87929	109.19094	B 1	40	0.00000
1246	122 PHE	CE1	-32.49275	-35.34303	110.14035	D1	40	0.00000
1247	122 PHE	CE2	-34,60187	-36.52254	110.55431	B1	40	0.00000
1248	122 PHE	CZ	-33.51285	-35.75823 -38.62424	105 60697	Bl	40	0.00000
1249	122 PHE	C	-34.3/844	-38.79208	104 65587	Bl	40	0.00000
1250	122 PHE 123 ASP	0 N	-31 06201	-39.60796	106.15389	Bl	41	0.00000
1251 1252	123 ASP	н	-31.23184	-39.47015		B 1	41	0.00000
1253	123 ASP	CA	-32.08552	-40.93522	105.58825	Bl	41	0.00000
1254	123 ASP	СВ	-30.85171	-41.81726	105.76445	Bl	41	0.00000
1255	123 ASP	CG	-29,93161	-41.56041	104.59405	Bl	41	0.00000
1256	123 ASP	ODl	-28.81173	-41.11310	104.81757	Bl	41	0.00000
1257	123 ASP	OD2	-30.34905	-41.80191	103.45856	Bl	41	0.00000
1258	123 A9P	С	-33.32362	-41.6361B	106.09965	B1	41 41	0.00000
1259	123 ASP	0	-34.00966	-41.21121	107.02319	BI	42	0.00000
1260	124 SER	И	-33.62443	-42.74250	105.41429	BJ DI	42	0.00000
1261	124 SER	Н	-32.95329	-43.10971 -43.35498	104.78900	B)	42	0.00000
1262	124 5ER	CY	-34.94423	-44.37028	104 46149	81	42	0.00000
1263	124 SER	CB	-35.18//9	-44.70776	104.37671	Bl	42	0.00000
1264	124 SER	OG	-36.3/343	-44.87590	105.27413	Bl	42	0.00000
1265 1266	124 SER 124 SER	,HG	-35.31660	-44.01057	106.93634	Bl	42	0.00000
1267	124 SER	0	-36.33538	-44.42223	107.22372	B1	42	0.00000 -
1268	125 ASP	N	-34.16447	-44.10325	107.74330	Bl	43	0.0000
1269	125 ASP	н	-33.26228	-43.73425	107.50494	81	43	0.00000
1270	125 ASP	CA	-34.30492	-44.60689	109.10471	B 1	43	0.00000
1271	125 ASP	CB	-32.96210	-45.24580	105.50620	Bl	43	0.00000
1272	125 ASP	CG	-31.82155	-44.23245	105.55779	B :	43	0.00000
1273	125 ASP	021	-31.89194	-43.19553	108.89075	51	43	0.00000 0.00000
1274	125 ASP	OD2	-30.87249	-44.45676	110.29614	P1	43	0.00000
1275	125 ASP	С	-34.65374	-43.52244	110.10428	57	43 43	c.00000
1276	125 ASP	0	-35.10540	-43.78441	:11.22872	21	• 3	0.0000

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1277	126	VAL	ы	- 51615	5 -42.26751	109.64790	B1	44	0.0000
1278		VAL	н	2093ء دن-	-42.12308	108.70456	Bl	4 4	0.00000
1279	126	VAL	CY	-34.59346	5 -41.07914	120.50496	Bl	4.4	0.00000
1280		VAL	CB	-36.04814	-40.82462	110.97592	Bl	4.4	0.00000
1281		VAL	CG1	-36.22164	-39.41758	111.54094	Bl	44	0.00000 0.00000
1282		VAL	CG2	-37.04494	-41.00281	109.82481	B1	4 4 4 4	0.00000
1283		VAL	C	-33.5762	5 -41.13739 -40.75401	111.01701	191	44	0.00000
1284		AYT	0	-33.78813	-41.66443	111.25324	Bl	45	0.0000
1285, 1286	127	GLY	H	-32.41347	-42.01122	110.31755	Bl	45	0.00000
1287		GLY	CA	-31 32244	-41.88201	112.19529	B1	4.5	0.00000
1288		GLY	c	-29.94594	-41.56889	111.62542	B1	45	0.00000
1289		GLY	ō	-29:03055	-41.21776	112.35962	Bl	45	0.00000
1290		GLU	N	-29.81143	-41.65704	110.29545	Bl	4 6	0.00000
1291	128	GLU	H	-30.48653	-42.13586	109.72534	B1	46	0.00000
1292	128	GLU	СX	-28.56256	-41.14297	109.73120	B1	4 6 4 6	0.00000 0.00000
1293	128	GLU	CB		-42.29481		ומ מ	46	0.00000
1294	128	GLU	CG		-42.25781 -43.28422			46	0.00000
1295		GLU	CD	-25.44161 -25.44161	-42.88073	108.57033	B1	46	0.00000
1296	128 · 128	GLU	OE2	-75 71010	-44.47773	109 00348	Bl	46	0.00000
1297 1290	128	GLU	C		-39.98012			46	0.00000
1299		GLU	0		-39.57375		Bl	46	0.00000
1300	129	TYR	N		-39.42619			47	0.00000
1300	129	TYR	H		-39.83420			47	0.00000
1302	129	TYR	CA.	-27.50019	-38.29454	107.46434	B1	47	0.00000
1303		TYR	CB		-37.15326		21	47	0.00000
1304	129	TYR	CG	-27.30857	-36.34222	109.09443	Bl	47	0.0000
1305		TYR	CD1	-26.67181	-36.21958	110.34811	Bl	47	0.00000
1306	129	TYR	CE1	-27.25566	-35.42131	111.35148	Bl	47	0.00000
1307		TYR	CD2		-35.67832		Bl	47	0.00000
1308		TYR	CE2		-34.88063	109.83571		. 47	0.00000
1309		TYR	CZ		-34.75726	111.08860		47	0.00000
1310		TYR	OH		-33.98459	•		47	0.00000
1311		TYR	нн	-29.70394	-33.40059 -38.64304			47	0.00000
1312 1313		IYR IYR	0	-25.62531	-39.04407	105.10492	B1	47	0.00000
1314		ARG	и		-38.38162	105.08581		4 B	0.00000
1315		ARG	н		-38.10451	105.16735	Bl	48	0.00000
1316		ARG	CA		-38.41254			48	0.00000
1317		J RG	СВ	-27.85650	-38.84699	102.71647		48	0.00000
1318	130	arg	CG	_ ,,	-39.00526	101.34112		48	0.00000
1319		NRG	CD		-39.41974	100.29808	Bl	48	0.00000
1320		ARG	NE		-39.38333	98.95629		4 B 4 B	0.00000 0.00000
1321	130		HE		-38.85034	98.82833		48	0.00000
1322 1323	130 . 130		CZ		-39.96348 -39.81313	96.69917	B1	48	0.00000
1323	130	-	NH1 HH11		-40.23513	,95.90996		48	0.00000
1325	130				-39.25297		Bl	48	0.00000
1326		ARG	NH2		-40.68730	98.14392	Bl	48	0.00000
1327	130			-29.90446		97.39118	Bl	48	0.00000
1328	130			-29.72741	-40.81930	99.07957	Bl	48	0.00000
1329	130	ARG	С		-37.06053	103.45986		48	0.00000
1330	130		0		-36.11772	103.07688	B1	48	0.00000
1331	131		ĸ	-24.95816	-36.99899	103.58668		49	0.00000 0.00000
1332	131		H	-24.45069	-37.77908	103.95192	D.T.	49 49	0.00000
1333	131		CA	-24.28607	-35.84894	102.98902		49	0.00000
1334 1335		ALA	CB C		-35.44271 -36.20633	103.80801		49	0.00000
1335	131 /	ALA =1.1	0	-23.17892		101.36065		49	0.00000
1336	132 1		ĸ	-24.28495		100.63154	Bl	50	0.00000
:338	132		16	-24.78984		100.85567	Bl	50	0.0000
1339	132		CA	-24.05930	-35.79619	99.24192	Вl	50	0.00000
1340	132 1		CB	-25.12578		98.36560	B 1	5 C	0.00000

./DR1_	MIKŽ . CPD		تعد 145	25 14:58:4	8 1993	22	
1341	132 VAL	CG1	5.02505	-35.45687	96.87606	22 30	.D. 570 0300
1342	132 VAL	CG2	-26.51048	-35.48853	98.85777	B1 50	0.00000
2243	135 ለሃን	C	-22.64446	-35.53393	98.73419	B1 50	0.00000
1344	132 VAI	0	-22.12598	-36.19257	97.84114 99.35299		0.00000
1345	133 THR	N	-22.01646	-34.53567 -34.04272	100 11190		0.00000
1346 1347	133 THR 133 THR	CY H		-34.13851	98.91128		0.00000
1348	133 THR	CB	-20.84979	-33.09962	97.76600		0.00000
1349	133 THR	061	-19.58519	-32.56706	97.34397		0.00000
1350	133 THR	HG1	-19.69018	-32.10136	96.50551		0.00000
1351	133 THR	CG2	-21.81730	-31.96947	98.12788		0.00000 0.00000
1352	133 THR	С	-19.91735	-33.59750	100.10846	B1 51 B1 51	0.00000
1353	133 THR	0	-20.48697	-33.30952 -33.44216	99.91793	B1 52	0.00000
1354	134 GLU 134 GLU	И Н.	-18.00113	-33.65725	99,02398		0.00000
1355 1356	134 GLU	CA	-17.75238	-32.88738	100.97647		0.00000
1357	134 GLU	CB	-16.30452	-32.81284	100.49482	B1 52	0.00000
1356	134 GLU	CG	-15.76229	-34.17073	100.03844	B1 52	0.00000
1359	134 GLU	CD	-14.31377	-34.04410	99.60488		0.00000
1360			13.54341	-34.96158	99.88007	B1 52 B1 52	0.00000
1361	134 GLU	OE2	-13.96021	-33.03519	98.99487	-	0.00000
1362	134 GLU	C	-18.18601	-31.51310 -31.11916	101.43728		0.00000
1363 1364	134 GLU 135 LEU	о и	-18 87611	-30.80674	100.55812		0.0000
1365	135 LEU	B	-18.92762	-31.13340	99.61398		0.00000
1366	135 LEU	CA.	-19.55727		100.94931	B1 53	0.0000
1367	135 LEU	CB	-20.29914	-29.06218	99.70959	B1 53	0.00000
1368	135 LEU	CG	-20.12211	-27.57937	99.38843		0.00000
1369	135 LEU	CD1	-20.93390	-26.68772	100.32666	B1 53	0.00000 0.00000
1370	135 LEU	CD2	-18.63204		.99.36436		0.00000
1371	135 LEU	С	-20.53099	-29.72915	102.11412	B1 53	0.00000
1372	135 LEU	0	-20.60025	-28.93514 -30.82221	102.04403		0.00000
1373 1374	136 GLY 136 GLY	N : H .	-21.14062	-31,49656	101.30480	B1 54	0.0000
1375	136 GLY	CA	-22.25373	-31.10000	103.08935	B1 54	0.00000
1376	136 GLY	C	-21.66227	-31.85099	104.26834	B1 54	0.00000
1377	136 GLY	0	-22.17280	-31.82013	105.38053	B1 54	0.00000
1378	137 ARG	N	-20.54192	-32.53644	104.01306	B1 55 B1 55	0.00000 0.00000
1379	137 ARG	H	-20,17033	-32.57514	103.00290		0.00000
1380	137 ARG 137 ARG	CA	-19.94165	-33.32437 -34.02762	104.58318	B1 55	0.00000
1381 1382	137 ARG 137 ARG	CB CG	-18 13008	-35.07893	105.55025		0.0000
1383	137 ARG	CD	-19.16855	-36.15719	105.85371	B 1 55	0.00000
1384	137 ARG	NE.	-18.66781	-37.06640	106.87657	B1 55	0.00000
1385	137 ARG	HE	-17.93002	-36.71980	107.47204	B1 55	0.00000
1386	137 ARG	CZ	-19.21896	-38.26612	107.07086	B1 55	0.00000 0.00000
1387	137 ARG	NH1	-18.67934	-39.07643	107.97489	B1 55 B1 55	0.00000
1388	137 ARG		-19.03969	-39.99255 -38.75541	108.14749		0.00000
1389 1390	137 ARG 137 ARG	HH12 NH2	-1/.88/63	-38.64369	106.37214		0.00000
1391	137 ARG	HH21	-20.72192	-39.53705	106.49834	B1 55	0.0000
1392	137 ARG	HH22	-20.68879	-38.01490	105.70437	B1 55	0.00000
1393	137 ARG	c,	-19.76338	-32.65026	106.46929	B1 55	. 0.00000
1394	137 ARG	0	-20.33202	-33.13638	107.44172	B1 55 B1 56	0.00000
1395	138 PRO	1 7	-19.03095	-31.51331	105.562//		0.00000
1396	138 PRO	CD	-18.26903	-30.78028 -30.86885	103.33102		0.00000
1397 1398	138 PRO 138 PRO	CA CB	-18 00866	-29.66775	107.61418	51 56	0.00000
1398	138 PRO	CG	-17.22921	-30.01427	106.35389		0.00000
1400	138 PRO	c	-20.24853	-30.39451	108.45452	er 55	0.00000
1401	138 PRO	ō	-20.38873	-30.20105	109.65243	31 56	0.00000
1402	139 ASP	N	-21.22822	-30.19487	107.56513	31 57	0.00000
1403	139 ASR	H	-21.12955	-30.45258	106.60295	91 57 91 57	0.00000 C.00000
1404	139 ASP	CA	-22.54445	-29.76374	198.07035	2:)·	3.00000

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1405	139 ASP	СВ	3 2048	-29 2279	4 106.80729 E	37.	~0~0 ∞ 00₫
1406	139 ASP		-24.5068	0 -28.4415	0 107.23406	1 57	0.00000
1.07	139 ASP	OD1	-24.33925	5 -27.2595	5 107.53698 E	13 57	0.00000
1408	139 ASP	002	-25.5959	0 -29.0093	7 107.24891 E	13 57	0.00000
1409	139 ASP	C	23.2900	9 -30.9065	1 108.70646 E	57	0.00000
1410	139 ASP		-23.84420	3 -30.7994	4 109.79628 E	11 57 11 58	0.00000 0.00000
1411	140 ALA		-23.1886	7 -32.0093	0 108.04975 E 5 107.13514 E	1 58	0.00000
1412	140 ALA 140 ALA	H CA	-22.7774	-32.0934	3 108.70194 2		0.00000
1413	140 ALA		-23.0414	-34.5128	5 107.80148 E	1 58	0.0000
1415	140 ALA	c	-22.96994	-33.5259	1 110.04660 E	1 58	0.00000
1416	140 ALA	ō	~23.61501	-33.62460	0 111.08086 E	1 58	0.00000
1417	141 GLU	14	-21.63141	-33.53710) 110.01537 E	1 59	0.00000
1418	141 GLU	H	-21.14260	-33.44422	109.14251 B	1 59	0.00000
1419	141 GLU	CA	-20.88133	-33.73045	111.26272 B	1 59	0.00000 0.00000
1420	141 GLU	CB	-19.38545	-33.75474	110.92637 B 109.97183 B	1 59 1 59	0.00000
1421	141 GLU	CG	-19.08157	-34.92133	109.97103 B	1 59	0.00000
1422	141 GLU 141 GLU	CD OE1	-17.05005	-34.91070	108.96182 B	1 59	0.00000
1424	141 GLU	· OE2	·· -16.99658	-33.87375	-109.50288 B	1 59	0.00000
1425	141 GLU	c	-21.20315	-32.72395	112.36696 B	1 59	0.00000
1426	141 GLU	ŏ	-21.35204	-33.05368	113.53869 B	າ 59	0.0000
1427	142 TYR	N	-21,39109	31.47026	.111.93865 B	1 60	0.00000
1428	142 TYR	н	-21.16858	-31.22939	110.99159 B	1 60	0.00000
1429	142 TYR	CA	-21.91640	-30.42572	112.82625 B	1 60	0.00000
1430	142 TYR	CB	-22.17510	-29.16770	111.95478 B	1 60 1 60	0.00000 0.00000
1431	142 TYR	CG	-22.15441	-27.86866	112.68902 B 112.88286 B	1 60	0.00000
1432 1433	142 TYR 142 TYR	CD1 CE1	-20.91930	-21.21121	113.49633 B	1 60	0.00000
1433	142 TYR	CD5	-23.36373	-27.27306	113.11310 B	1 60	0.00000
1435	142 TYR	CE2	-23.33600	-26.00211	113.72688 B	1 60	0.00000
1436	142 TYR	CZ	-22,10013	-25.34007	113.91274 B	1 60	0.00000
1437	142 TYR	ОН	-22.06472	-24.08718	114.49226 B	1 60	0.00000
1438	142 TYR	HH	-22.95958	-23.75980	114.62492 B	1 60	0.00000
1439	142 TYR	Ċ	-23.20365	-30.84932	113.52485 B	1 60 1 60	0.00000
1440	142 TYR	0	-23.33185	-30.86333	114.74427 B 112.69102 B	1 61	0.00000
1441	143 TRR 143 TRP	h n	-24.16819	-31.24330	111.69669 B	1 61	0.00000
1443	143 TRP	CA	-25 46084	-31.65772	113.24427 B	61	0.00000
2444	143 TRP	CB	-26.46502	-31.82534	112.10045 B	61	0.0000
1445	143 TRP	CG	-26.82927	-30.51319	111,43167 B	1 61	0.00000
1446	143 TRP	CD2	-27.59514	-30.35383	110.26190 B	1 61	0.00000
1447	143 TRP	CE2.	-27.68725	-28.88192	110.01923 B	1 61 1 61	0.00000
1448	143 TRP	CE3	-28.23171	-31.24445	109.37526 B	61	0.00000
1449 1450	143 TRP	CD1 NE1	-26.49404	-29.20400	111.01939 B	61	0.00000
1451	143 TRP	HE1	-26.86335	-27.27405	111.09738 B	61	0.00000
1452	143 TRP	CZ2	-28.41151	-28.41903	108.90296 B	61	0.00000
1453	143 TRP	CZ3	-28.94655	-30.73482	108.27096 B	1 61	0.00000
1454	143 TRP	CH2	-29.03488	-29.34388	108.03833 B	61	0.00000
1455	143 TRP	C·	-25.40824	-32.93379	114.07770 B	61	0.00000
1456	143 TRP	Ο,	-26.13451	-33.11650	115.04995 B	61	0.00000
1457	144 A6N	N	-24.46546	-33.80055	113.69236 B	L 62 L 62	0.00000
1458	144 ASN	X.	-23.94027	-33.62305	112.85783 E	62	0.00000
1459 1460	144 ASN 144 ASN	CA CB	-24.10067	-35.35UUU	113.73882 B		0.00000
1451	144 ASN	CG	-23.20030 -23 89541	-36.68861	112.61740 B	62	0.00000
1462	144 ASN	OD1	-23.68155	-36.47075	111.43371 B	62	0.00000
1463	144 ASN	ND2	-24.72776	-37.64097	113.02365 B	62	0.00000
1464	144 ASN	HD21	-24.89338	-37.61926	113.99276 B	62	6.00000
1465	144 ASN		-25.19584	-39.20406	112.34455 B	62	. C. 00000 0. 00000
1466	144 ASN	C	-23.49875	-34.69497	115.82591 B1	62	o.00000
1467	144 ASN	l: 0	-23.43003	-33.34034	115.97257 B	63	0.00000
	145 SER		-44.57004	- 22.45040			

FIG. 24

.,	يان ، - درندي	ن	شه د ددد	:	دبده ٥٠	24	
1469	145 821	ч.	22.9783	0 -32.8042	8 115.21961 81	69	· •0 :-0 0 -0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0
1470	145 SZF		22.3800	4 -33.1462	1 117.25865 B1	63	0.00000
3477	145 SEF	R CB	-20.9192	1 -32.7305	4 117.01176 B1	63	0.00000
1472	145 SEF	R OG	-20.1898	3 -32.6418	0 118.24511 B1	63	0.00000
1473					5 118.91065 B1	63	0.00000
1474		₹ C			4 118.06161 B1	63	.0.0000
1475			-22.6740	1 -31.6828	7 119.13941 B1	63	0.00000
1476					7 117.51252 Bl	64	0.00000
1477					4 116.64774 B1	64	0.00000
2478	146 GLN				6 118.19522 Bl	64	0.00000
1479	146 GLN		-24.8631	5 -29.2930	1 117.23892 B1	64	0.00000
1480	146 GLN				5 117.72882 B1	64	0.00000
1481	146 GLN				2 117.25234 B1	64	0.00000
1482					9 118.00911 B1 8 115.94329 B1	64 64	0.00000
1483 1484	146 GLN 146 GLN				9 115.33276 B1	64	0.00000
1485	146 GLN	HE21	-20.3000:	7 -20.2034; 7 -27 8899°	7 115.56134 B1	64	0.00000
1486	146 GLN				4 118.57642 B1	64	0.00000
1487	146 GLN				9 117.77294 B1	64	0.00000
1488	147LYS				6 119.82371 B1	65	0.00000
1489	147 LYS				120.47663 B1	65	0.00000
1490	147 LYS				120.21964 B1	65	0.00000
1491	147 LYS	CB .	-28.27258	32.00853	121.37895 B1	65	0.00000
1492	147 LYS	CG ·	-29.34716	-33.02726	121.84203 B1	65	0.00000
1493	147 LYS	CD :	-28.88674	-34.49172	2 121.97244 B1	£5	0.00000
1494	147 LYS	CE .	-29.96618	-35.46785	121.47444 B1	65	0.00000
1495	147 LYS	NZ ·	-29,40221	-36.81034	121.26907 Bl	€5	0.00000
1496	147 LYS	H21 -	-30.04886	-37.42223	120.71767 B1	65	0.00000
1497	147 LYS				120.68922 B1	65	0.0000
1498	147 LYS	H23 -	-29.15199	-37.28892	122.15073 B1	65	0.0000
1499	147 LYS				119.09340 Bl	65	0.00000
1500	147 LYS				118.83883 Bl	65	0.00000
1501	148 ASP		•		118.45822 B1	66	0.00000
1502	148 ASP				118.75300 Bl	66	0.00000
1503	148 ASP				117.48362 Bl	66 66	0.00000
1504	148 ASP	CB ~	27.88927	-36.37833	118.17059 B1 119.21585 B1	66	0.00000
1505 1506	148 ASR 148 ASP	CG -	25 60280	-36.37330	119.05193 B1	66	0.00000
1507	148 ASP	OD2 -	27.07650	-37.20742	120.23524 B1	66	0.00000
1508	148 ASP	c	29.08887	-34.63043	116.79645 B1	66	0.00000
1509	148 ASP	0 -	30.17136	-35.14808	117.04951 B1	€6	0.00000
1510	149 LEU				115.93941 B1	67	0.00000
1511	149 LEU			-33.30949		67	0.00000
1512	149 LEU	CA -	30.07076	-32.84836	115.45008 B1	67	0.00000
1513	149 <u>L</u> EŲ	СБ -	29.45399	-31.74267	114.57360 B1	67	0.00000
1514	149 LEU				114.02475 B1	67	0.00000
1515	149 LEU				112.65290 B1	67	0.00000
1516	149 LEU		31.34761	-30.09615	115.02072 B1	67	0.00000
1517	149 LEU				114.80952 B1	67	0.00000
1518	149 LEU	0 -	32,36472	-33.37859	114.83807 Bl	67	0.00000 0.00000
1519	150 LEU				114.32363 Bl	68 68	0.00000
1520	150 LEU	K -	29.75579	-34.94139	114.11621 B1	68	0.00000
1521 1522	150 LEU 150 LEU	CA -	30 34340	-30.02822	114.12850 B1 114.15286 B1	68	0.00000
1522	150 LEU				114.15266 B1 112.89569 B1	68	0.00000
1524	150 LEU				112.89369 B1 113.18796 B1	68	0.00000
1525	150 LEU				111.73915 B1	68	0.00000
1526	150 LEU	C -:	32.74973	-36,17247	115.10785 B1	65	0.00000
1527	150 LEU				114.70350 B1	63	0.00000
1528	151 GLU				116.39428 B1	6.5	0.00000
1529	151 GLU	н - ;	31.52943	-36.57284	116.72119 B1	6 è	0.00000
1530	151 GLU	CA -3	33.61295	-36.59512	117.30950 E1	6.9	0.00000
1331	151 GLU	CB3	33.19489	-36.98331	118.72928 81	65	0.00000
1532	721 CTA	CG -:	32.69081	-36.41906	118.86324 91	6 ?	0.00000

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. / 🍱	CERTHS CE	B	மோ உடி	25 14:58	:48 1993	25	
253	27 151 00						
. 131		U CD	-31.1928	37 -38.444	64 118.7049	7 El 69	
153		U OE1			29 117.5814		· -
				22 -38.564		9 Bl 69	
153			-34.5506	7 -35.411	68 117.4385	7 Bl 69	0.00000
153		-		0 -35.573		7 B1 69	0.00000
153			-33.9860	1 -34.197	80 117.3558	5 Bl 70	0.00000
153			-33.0032	1 - 34.102	56 117.18602	B1 70	0.00000
154			-34.8965	6 -33.047	52 117.40695	B1 70	0.0000
154		N CB	-34.1568	0 ~31.716	46 117.32018	B1 70	0.00000
154	2 152 GL	N CG	-33.4615	9 -31.269	00 118.60094	B1 70	0.00000
154	3 152 GL	N CD	-33.2337	6 -29.774	24 118.49387	B1 70	0.00000
154	4 152 GL	N OE1	-32.1306	8 -29.270	19 118.36008	B1 70	0.00000
154		NE2			52 118.54711		0.00000
154	6 152 GL	N HE21	-35.2460	7 -29.465	11 118.65663	B1 70	0.00000
154	7 152 GL				8 118.46911		0.00000
154	9 152 GL	1 C			6 116.27182		0.00000
154	9 152 GLA	1 0			1 116.43607		0.00000
1550	153 ARG	N			9 115.08552		0.00000
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FIG. 30

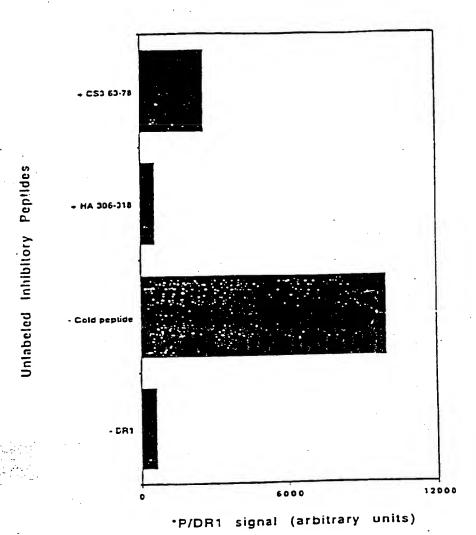
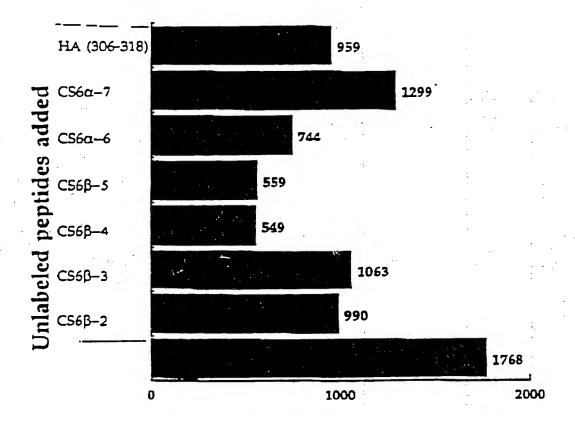


FIG. 31

30/31

Inhibition of 125 I HA (306-318)/DRI by unlabeled CSG of and B pertides



*HA/DR1 compact dimer signal (densitometric units)

INTERNATIONAL SEARCH REPORT

International application No. PCT/US94/05697

A. CLASSIFICATION OF SUBJECT MATTER IPC(5) :A61K 39/00, 39/02, 39/12, 37/02, 35/14 US CL :424/185.1, 186.1, 190.1, 242.1; 530/327, 326, 333, 334, 388.75			
According to International Patent Classification (IPC) or to both national classification and IPC B. FIELDS SEARCHED			
Minimum documentation searched (classification system followed by classification symbols)			
U.S. : 424/185.1, 186.1, 190.1, 242.1; 530/327, 326, 333, 334, 388.75			
Documentation searched other than minimum documentation to th	e extent that such documents are included in the fields searched		
Electronic data base consulted during the international search (name of data base and, where practicable, search terms used)			
C. DOCUMENTS CONSIDERED TO BE RELEVANT			
Category* Citation of document, with indication, where a	ppropriate, of the relevant passages Relevant to claim No.		
issued 15 April 1993, Nauss et al Peptides in a Structural Homology	The Journal of Immunology, Volume 150, No. 8, Part II, issued 15 April 1993, Nauss et al., "Binding Interactions of Peptides in a Structural Homology Model of the DR1 Class MHC ", page 41A, Abstract 221, see entire abstract.		
X Nature, Volume 358, issued 27 A "Predominant Naturally Processed DR1 are derived from MHC-re Heterogenous in Size", pages 764 2, and Table 3.	Peptides Bound to HLA- lated Molecule and are 1, 3-7		
·			
X Further documents are listed in the continuation of Box C. See patent family annex.			
Special categories of cited documents: "T" later document published after the international filling date or priority date and sot in conflict with the application but cited to understand the principle or theory underlying the invention.			
to be of particular relevance	principle or theory underlying the invention "X" document of particular relevance; the claimed invention cannot be		
*E' earlier document published on or after the international filing date *L' document which may throw doubts on priority claim(s) or which is	considered novel or cannot be considered to involve an investive step when the document is taken alone		
cited to establish the publication date of another citation or other special reason (as specified)	"Y" document of particular relevance; the claimed invention cannot be		
O document referring to an oral disclosure, use, exhibition or other means being obvious to a person skilled in the art			
P document published prior to the international filing date but later than the priority date claimed	'P' document published prior to the international filing date but later than "g" document member of the same patent family		
Date of the actual completion of the international search Date of mailing of the international search report			
01 SEPTEMBER 1994 1 3 SEP 1994.			
Name and mailing address f the ISA/US Commissioner of Patents and Trademarks Authorized fficer			
Commissioner of Patents and Trademarks Box PCT Washington, D.C. 20231 H. Sidberry			
Facsimile No. (703) 305-3230	Telephone No. (703) 308-0196		

INTERNATIONAL SEARCH REPORT

International application No. PCT/US94/05697

Category*	Citation f document, with indication, where a	ppropriate, of the relevant passages	Relevant to claim No
Y	The Journal of Immunology, Volume 150, No. 2, issued 15 January 1993, Boehncke et al., "The Importance of Dominant Negative Effects of Amino Acid Side Chain Substitution in Peptide-MHC Molecule Interactions and T Cell Recognition", pages 331-341, see Abstract, on page 331.		8-11
ζ	The EMBO Journal, Volume 9, No. 6 al., "Peptide binding to HLA-DR1: a substituted to alanine retains MHC bir page 1798, page 1800, figure 4, and p	Peptide with most residues adding", pages 1797-1803, see	512
?	Nature, Volume 332, issued 28 April hypothetical model of the foreign antiphistocompatibility molecules", pages 8	gen binding site of Class II	1, 3, 4
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	Fig. 1	•	
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INTERNATIONAL SEARCH REPORT

International application No. PCT/US94/05697

Box 1 Observations where certain claims w re f und unsearchable (Continuation f item 1 f first sheet)
This international report has not been established in respect of certain claims under Article 17(2)(a) for the following reasons:
Claims Nos.: 2 because they relate to subject matter not required to be searched by this Authority, namely:
Claim 2 is directed to a computerized model which encompasses scientific theory and computer programs to the extent that the International Searching Authority is not equipped to search prior art concerning such programs. Accordingly claim 2 is withdrawn from search under PCT Rule 39 and PCT Article 17(2)(a)(i).
Claims Nos.: because they relate to parts of the international application that do not comply with the prescribed requirements to such an extent that no meaningful international search can be carried out, specifically:
Claims Nos.: because they are dependent claims and are not drafted in accordance with the second and third sentences of Rule 6.4(a).
Box 11 Observations where unity of invention is lacking (Continuation of item 2 of first sheet)
This International Searching Authority found multiple inventions in this international application, as follows:
1. As all required additional search fees were timely paid by the applicant, this international search report covers all searchable claims.
2. As all searchable claims could be searched without effort justifying an additional fee, this Authority did not invite payment of any additional fee.
3. As only some of the required additional search fees were timely paid by the applicant, this international search report covers only those claims for which fees were paid, specifically claims Nos.:
<u> </u>
4. No required additional search fees were timely paid by the applicant. Consequently, this international search report is restricted to the invention first mentioned in the claims; it is covered by claims Nos.:
Remark on Protest
N protest accompanied the payment f additional search fees.

Form PCT/ISA/210 (continuation of first sheet(1))(July 1992)*